

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081821\
 Data File : BP006748.D
 Acq On : 18 Aug 2021 16:57
 Operator : CG/JU
 Sample : M3394-04MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SWCS-01-0205MSD

Manual Integrations
 APPROVED

mohammad

8/19/2021 1:27:35 PM

Quant Time: Aug 18 18:23:26 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 18:17:59 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	164489	20.000	ng	0.00
21) Naphthalene-d8	10.487	136	669808	20.000	ng	# 0.00
39) Acenaphthene-d10	14.345	164	374821	20.000	ng	0.00
64) Phenanthrene-d10	17.104	188	755975	20.000	ng	# 0.00
76) Chrysene-d12	21.216	240	737837	20.000	ng	# 0.00
86) Perylene-d12	23.527	264	781915	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	868041	81.055	ng	0.00
7) Phenol-d6	6.893	99	1107733	72.816	ng	0.00
23) Nitrobenzene-d5	8.863	82	760951	52.438	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	344567	87.961	ng	0.00
45) 2-Fluorobiphenyl	12.969	172	1369696	54.672	ng	0.00
79) Terphenyl-d14	19.686	244	2079803	56.482	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.246	88	196310	42.105	ng	95
3) Pyridine	3.634	79	488643	35.658	ng	97
4) n-Nitrosodimethylamine	3.552	42	189276	36.595	ng	# 73
6) Aniline	7.040	93	526642	31.870	ng	97
8) 2-Chlorophenol	7.269	128	512906	44.264	ng	94
9) Benzaldehyde	6.852	77	365686	39.939	ng	93
10) Phenol	6.916	94	637799	41.811	ng	99
11) bis(2-Chloroethyl)ether	7.140	93	482724	41.676	ng	93
12) 1,3-Dichlorobenzene	7.587	146	547530	45.338	ng	97
13) 1,4-Dichlorobenzene	7.740	146	557785	45.509	ng	99
14) 1,2-Dichlorobenzene	8.046	146	529510	45.023	ng	97
15) Benzyl Alcohol	7.952	79	471534	41.332	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.240	45	531824	38.585	ng	96
17) 2-Methylphenol	8.157	107	421042	43.712	ng	99
18) Hexachloroethane	8.769	117	220757	45.505	ng	92
19) n-Nitroso-di-n-propyla...	8.516	70	397168	38.914	ng	94
20) 3+4-Methylphenols	8.487	107	563349	42.969	ng	97
22) Acetophenone	8.528	105	771929	43.457	ng	# 98
24) Nitrobenzene	8.904	77	610247	40.455	ng	92
25) Isophorone	9.434	82	1031738	39.584	ng	95
26) 2-Nitrophenol	9.610	139	263913	48.074	ng	93
27) 2,4-Dimethylphenol	9.681	122	444877	48.394	ng	97
28) bis(2-Chloroethoxy)met...	9.916	93	616049	44.175	ng	99
29) 2,4-Dichlorophenol	10.140	162	423509	45.504	ng	93
30) 1,2,4-Trichlorobenzene	10.351	180	450467	45.042	ng	97
31) Naphthalene	10.534	128	1542031	42.844	ng	100
32) Benzoic acid	9.846	122	346635	48.742	ng	91
33) 4-Chloroaniline	10.657	127	259126	17.791	ng	96
34) Hexachlorobutadiene	10.816	225	268574	46.085	ng	98
35) Caprolactam	11.469	113	152360m	45.400	ng	
36) 4-Chloro-3-methylphenol	11.798	107	520982	44.058	ng	93
37) 2-Methylnaphthalene	12.157	142	1056360	43.313	ng	99
38) 1-Methylnaphthalene	12.375	142	970003	41.852	ng	98
40) 1,2,4,5-Tetrachloroben...	12.528	216	457495	46.635	ng	# 95
41) Hexachlorocyclopentadiene	12.498	237	335829	64.545	ng	98
43) 2,4,6-Trichlorophenol	12.775	196	316854	47.302	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081821\
 Data File : BP006748.D
 Acq On : 18 Aug 2021 16:57
 Operator : CG/JU
 Sample : M3394-04MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SWCS-01-0205MSD

Manual Integrations
 APPROVED

mohammad

8/19/2021 1:27:35 PM

Quant Time: Aug 18 18:23:26 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 18:17:59 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.845	196	345150	46.550	ng	# 90
46) 1,1'-Biphenyl	13.181	154	1215165	42.832	ng	97
47) 2-Chloronaphthalene	13.216	162	959697	43.927	ng	94
48) 2-Nitroaniline	13.434	65	346791	44.978	ng	87
49) Acenaphthylene	14.069	152	1595424	45.288	ng	99
50) Dimethylphthalate	13.822	163	1172012	42.956	ng	99
51) 2,6-Dinitrotoluene	13.939	165	259604	42.641	ng	# 81
52) Acenaphthene	14.410	154	925449	43.165	ng	98
53) 3-Nitroaniline	14.263	138	158876	23.539	ng	82
54) 2,4-Dinitrophenol	14.481	184	134494	42.244	ng	# 87
55) Dibenzofuran	14.745	168	1440372	42.711	ng	93
56) 4-Nitrophenol	14.592	139	547010	93.337	ng	89
57) 2,4-Dinitrotoluene	14.728	165	370883	45.424	ng	# 80
58) Fluorene	15.398	166	1185629	43.985	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.981	232	289999	46.671	ng	# 72
60) Diethylphthalate	15.186	149	1272873	44.418	ng	97
61) 4-Chlorophenyl-phenyle...	15.398	204	538249	44.094	ng	# 87
62) 4-Nitroaniline	15.434	138	267547m	37.357	ng	
63) Azobenzene	15.692	77	1433910	42.917	ng	92
65) 4,6-Dinitro-2-methylph...	15.492	198	92671	20.062	ng	73
66) n-Nitrosodiphenylamine	15.622	169	1001467	42.147	ng	99
67) 4-Bromophenyl-phenylether	16.298	248	326268	44.981	ng	# 90
68) Hexachlorobenzene	16.404	284	372931	45.180	ng	# 88
69) Atrazine	16.586	200	361245	48.720	ng	97
70) Pentachlorophenol	16.757	266	509107	97.209	ng	99
71) Phenanthrene	17.151	178	1849339	43.370	ng	99
72) Anthracene	17.239	178	1894976	45.990	ng	99
73) Carbazole	17.516	167	1766156	42.023	ng	98
74) Di-n-butylphthalate	18.086	149	2313735	48.321	ng	100
75) Fluoranthene	19.127	202	2110067	44.087	ng	95
77) Benzidine	19.322	184	1917823	103.873	ng	99
78) Pyrene	19.480	202	2170202	44.035	ng	99
80) Butylbenzylphthalate	20.374	149	1094667	50.512	ng	92
81) Benzo(a)anthracene	21.198	228	2079936	43.134	ng	99
82) 3,3'-Dichlorobenzidine	21.139	252	775912	61.293	ng	# 96
83) Chrysene	21.251	228	2014036	43.083	ng	99
84) Bis(2-ethylhexyl)phtha...	21.139	149	1632402	50.125	ng	# 97
85) Di-n-octyl phthalate	22.039	149	2856194	53.210	ng	# 97
87) Indeno(1,2,3-cd)pyrene	25.921	276	2603921	45.927	ng	# 91
88) Benzo(b)fluoranthene	22.827	252	2264603	44.563	ng	# 97
89) Benzo(k)fluoranthene	22.874	252	2065950	42.342	ng	# 97
90) Benzo(a)pyrene	23.427	252	2133913	46.816	ng	# 95
91) Dibenzo(a,h)anthracene	25.945	278	2218586	45.258	ng	# 95
92) Benzo(g,h,i)perylene	26.656	276	2173426	46.267	ng	# 92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081821\
Data File : BP006748.D
Acq On : 18 Aug 2021 16:57
Operator : CG/JU
Sample : M3394-04MSD
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SWCS-01-0205MSD

Manual Integrations
APPROVED

mohammad

8/19/2021 1:27:35 PM

Quant Time: Aug 18 18:23:26 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 16 18:17:59 2021
Response via : Initial Calibration

