

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
 Data File : BP006162.D
 Acq On : 09 Jul 2021 15:49
 Operator : CG/JU
 Sample : M2990-13MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-JMS

Manual Integrations
 APPROVED

mohammad
 7/12/2021 12:36:53 PM

Quant Time: Jul 09 18:35:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 07 16:08:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	184416	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	686073	20.000	ng	0.00
39) Acenaphthene-d10	14.510	164	338637	20.000	ng	0.00
64) Phenanthrene-d10	17.263	188	581393	20.000	ng	# 0.00
76) Chrysene-d12	21.345	240	506524	20.000	ng	# 0.00
86) Perylene-d12	23.745	264	621264	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	942764	87.962	ng	-0.01
7) Phenol-d6	7.058	99	1097608	77.423	ng	0.00
23) Nitrobenzene-d5	9.040	82	610656	56.222	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	294373	73.015	ng	0.00
45) 2-Fluorobiphenyl	13.134	172	1307296	57.071	ng	0.00
79) Terphenyl-d14	19.816	244	1755776	68.343	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.305	88	221964	52.782	ng	91
3) Pyridine	3.723	79	545636	51.065	ng	# 87
4) n-Nitrosodimethylamine	3.634	42	253845	60.400	ng	# 86
6) Aniline	7.205	93	817694	54.666	ng	98
8) 2-Chlorophenol	7.440	128	715776	57.022	ng	98
9) Benzaldehyde	7.016	77	342245m	44.233	ng	
10) Phenol	7.087	94	781062	57.365	ng	92
11) bis(2-Chloroethyl)ether	7.299	93	592043	57.544	ng	98
12) 1,3-Dichlorobenzene	7.758	146	768739	57.155	ng	96
13) 1,4-Dichlorobenzene	7.905	146	782420	56.959	ng	98
14) 1,2-Dichlorobenzene	8.222	146	741717	56.718	ng	99
15) Benzyl Alcohol	8.128	79	522395	60.951	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.399	45	664038	54.344	ng	90
17) 2-Methylphenol	8.334	107	523998	56.572	ng	94
18) Hexachloroethane	8.946	117	266303	58.234	ng	95
19) n-Nitroso-di-n-propyla...	8.687	70	394038	50.983	ng	98
20) 3+4-Methylphenols	8.663	107	694073	55.465	ng	92
22) Acetophenone	8.705	105	898452	57.536	ng	98
24) Nitrobenzene	9.087	77	607197	55.276	ng	91
25) Isophorone	9.616	82	1034506	50.172	ng	95
26) 2-Nitrophenol	9.799	139	359715	58.904	ng	# 84
27) 2,4-Dimethylphenol	9.863	122	573654	63.637	ng	96
28) bis(2-Chloroethoxy)met...	10.093	93	674098	58.993	ng	100
29) 2,4-Dichlorophenol	10.346	162	552871	55.819	ng	99
30) 1,2,4-Trichlorobenzene	10.540	180	589791	55.928	ng	97
31) Naphthalene	10.728	128	1950692	54.541	ng	99
32) Benzoic acid	10.069	122	375512	53.867	ng	93
33) 4-Chloroaniline	10.852	127	369074	26.733	ng	97
34) Hexachlorobutadiene	11.005	225	316926	54.583	ng	99
35) Caprolactam	11.669	113	167631	53.689	ng	# 85
36) 4-Chloro-3-methylphenol	11.993	107	557164	52.092	ng	98
37) 2-Methylnaphthalene	12.340	142	1311475	52.850	ng	97
38) 1-Methylnaphthalene	12.557	142	1195033	50.886	ng	98
40) 1,2,4,5-Tetrachloroben...	12.710	216	523344	62.302	ng	97
41) Hexachlorocyclopentadiene	12.681	237	352798	119.821	ng	98
43) 2,4,6-Trichlorophenol	12.957	196	359758	59.197	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
 Data File : BP006162.D
 Acq On : 09 Jul 2021 15:49
 Operator : CG/JU
 Sample : M2990-13MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-JMS

Manual Integrations
 APPROVED

mohammad
 7/12/2021 12:36:53 PM

Quant Time: Jul 09 18:35:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 07 16:08:13 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.040	196	382101	58.602	ng	# 94
46) 1,1'-Biphenyl	13.346	154	1463969	57.584	ng	98
47) 2-Chloronaphthalene	13.393	162	1137830	58.177	ng	98
48) 2-Nitroaniline	13.610	65	293664	60.424	ng	87
49) Acenaphthylene	14.234	152	1808686	57.431	ng	100
50) Dimethylphthalate	13.975	163	1312352	53.893	ng	99
51) 2,6-Dinitrotoluene	14.104	165	292538	53.286	ng	92
52) Acenaphthene	14.575	154	1042724	54.430	ng	98
53) 3-Nitroaniline	14.434	138	214353	38.538	ng	95
54) 2,4-Dinitrophenol	14.657	184	274711	93.560	ng	96
55) Dibenzofuran	14.910	168	1591248	53.358	ng	98
56) 4-Nitrophenol	14.775	139	461383	105.181	ng	# 82
57) 2,4-Dinitrotoluene	14.898	165	401868	54.351	ng	96
58) Fluorene	15.563	166	1273766	53.165	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.151	232	269393	52.803	ng	# 73
60) Diethylphthalate	15.334	149	1315398	52.673	ng	98
61) 4-Chlorophenyl-phenyle...	15.551	204	553181	52.668	ng	90
62) 4-Nitroaniline	15.604	138	276754	48.663	ng	# 84
63) Azobenzene	15.845	77	1091939	52.375	ng	96
65) 4,6-Dinitro-2-methylph...	15.669	198	178924	47.216	ng	90
66) n-Nitrosodiphenylamine	15.769	169	1077750	56.931	ng	100
67) 4-Bromophenyl-phenylether	16.451	248	321481	56.091	ng	93
68) Hexachlorobenzene	16.569	284	375946	55.261	ng	# 92
69) Atrazine	16.734	200	334501	56.282	ng	94
70) Pentachlorophenol	16.928	266	410725	130.810	ng	99
71) Phenanthrene	17.304	178	1833475	55.840	ng	100
72) Anthracene	17.398	178	1882679	58.529	ng	99
73) Carbazole	17.675	167	1741342	54.630	ng	99
74) Di-n-butylphthalate	18.216	149	2195397	57.220	ng	99
75) Fluoranthene	19.275	202	1956693	54.372	ng	95
77) Benzidine	19.457	184	1098442	73.133	ng	99
78) Pyrene	19.628	202	1994848	58.267	ng	98
80) Butylbenzylphthalate	20.486	149	998943	58.824	ng	94
81) Benzo(a)anthracene	21.333	228	1776472	55.452	ng	99
82) 3,3'-Dichlorobenzidine	21.263	252	590730	62.857	ng	# 97
83) Chrysene	21.386	228	1773185	56.007	ng	99
84) Bis(2-ethylhexyl)phtha...	21.245	149	1434480	57.229	ng	# 98
85) Di-n-octyl phthalate	22.157	149	2477566	56.121	ng	100
87) Indeno(1,2,3-cd)pyrene	26.274	276	2999788	65.493	ng	# 88
88) Benzo(b)fluoranthene	23.010	252	2047686	54.336	ng	# 95
89) Benzo(k)fluoranthene	23.063	252	2044701	55.510	ng	# 96
90) Benzo(a)pyrene	23.639	252	2112947	59.748	ng	# 94
91) Dibenzo(a,h)anthracene	26.286	278	2551729	64.699	ng	# 92
92) Benzo(g,h,i)perylene	27.051	276	2618850	67.470	ng	# 89

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

