

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011822\
 Data File : BP008841.D
 Acq On : 18 Jan 2022 16:04
 Operator : CG/JU
 Sample : N1154-01MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20220028-KEANSBURG-2-COMPMS

Quant Time: Jan 19 03:36:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 14 18:05:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	230860	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	864463	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	539417	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1154470	20.000	ng	0.00
76) Chrysene-d12	21.339	240	1222223	20.000	ng	0.00
86) Perylene-d12	23.756	264	1345896	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1832243	122.733	ng	0.00
7) Phenol-d6	7.028	99	2242734	124.060	ng	0.00
23) Nitrobenzene-d5	9.010	82	1301440	85.472	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	993079	126.478	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	2826179	69.093	ng	0.00
79) Terphenyl-d14	19.810	244	3839597	53.026	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.346	88	234492	38.807	ng	96
3) Pyridine	3.740	79	677441	53.529	ng	96
4) n-Nitrosodimethylamine	3.652	42	277162	46.679	ng	83
6) Aniline	7.175	93	568965	25.216	ng	97
8) 2-Chlorophenol	7.416	128	719965	45.282	ng	98
9) Benzaldehyde	6.987	77	380374	31.482	ng	97
10) Phenol	7.057	94	893993	48.507	ng	96
11) bis(2-Chloroethyl)ether	7.275	93	634794	43.290	ng	97
12) 1,3-Dichlorobenzene	7.740	146	799920	42.247	ng	98
13) 1,4-Dichlorobenzene	7.887	146	821420	42.805	ng	99
14) 1,2-Dichlorobenzene	8.204	146	781567	42.714	ng	99
15) Benzyl Alcohol	8.093	79	585015	46.128	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.381	45	698688	42.607	ng	99
17) 2-Methylphenol	8.304	107	567133	45.932	ng	96
18) Hexachloroethane	8.928	117	277036	42.535	ng	98
19) n-Nitroso-di-n-propyla...	8.663	70	465396	44.040	ng	94
20) 3+4-Methylphenols	8.634	107	756478	46.152	ng	96
22) Acetophenone	8.675	105	1008127	45.779	ng	# 96
24) Nitrobenzene	9.051	77	707242	45.983	ng	96
25) Isophorone	9.581	82	1240857	45.094	ng	98
26) 2-Nitrophenol	9.763	139	375206	46.403	ng	93
27) 2,4-Dimethylphenol	9.834	122	638107	46.285	ng	98
28) bis(2-Chloroethoxy)met...	10.063	93	783291	43.622	ng	99
29) 2,4-Dichlorophenol	10.304	162	690943	45.927	ng	99
30) 1,2,4-Trichlorobenzene	10.516	180	759930	43.122	ng	100
31) Naphthalene	10.704	128	2075656	44.208	ng	100
32) Benzoic acid	10.004	122	436129	55.123	ng	98
33) 4-Chloroaniline	10.816	127	318654	16.654	ng	99
34) Hexachlorobutadiene	10.992	225	462124	41.972	ng	98
35) Caprolactam	11.604	113	208725	50.328	ng	92
36) 4-Chloro-3-methylphenol	11.951	107	650514	47.929	ng	100
37) 2-Methylnaphthalene	12.316	142	1522153	44.347	ng	98
38) 1-Methylnaphthalene	12.534	142	1404204	44.571	ng	99
40) 1,2,4,5-Tetrachloroben...	12.686	216	865737	44.018	ng	99
41) Hexachlorocyclopentadiene	12.669	237	689550	68.518	ng	100
43) 2,4,6-Trichlorophenol	12.928	196	563329	46.325	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011822\
 Data File : BP008841.D
 Acq On : 18 Jan 2022 16:04
 Operator : CG/JU
 Sample : N1154-01MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20220028-KEANSBURG-2-COMPMS

Quant Time: Jan 19 03:36:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 14 18:05:00 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.004	196	609581	46.576	ng	97
46) 1,1'-Biphenyl	13.322	154	1976024	45.026	ng	100
47) 2-Chloronaphthalene	13.363	162	1511826	44.676	ng	99
48) 2-Nitroaniline	13.569	65	387850	51.619	ng	95
49) Acenaphthylene	14.210	152	2367261	46.302	ng	100
50) Dimethylphthalate	13.951	163	1841235	45.960	ng	100
51) 2,6-Dinitrotoluene	14.069	165	411223	48.399	ng	93
52) Acenaphthene	14.557	154	1400565	46.188	ng	99
53) 3-Nitroaniline	14.398	138	199792	23.910	ng	99
54) 2,4-Dinitrophenol	14.610	184	448071	130.396	ng	94
55) Dibenzofuran	14.892	168	2270546	45.952	ng	98
56) 4-Nitrophenol	14.722	139	727783	120.951	ng	92
57) 2,4-Dinitrotoluene	14.857	165	566764	51.860	ng	# 90
58) Fluorene	15.539	166	1856625	47.573	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.122	232	525167	48.712	ng	# 85
60) Diethylphthalate	15.316	149	1776734	46.741	ng	98
61) 4-Chlorophenyl-phenyle...	15.533	204	965162	45.398	ng	98
62) 4-Nitroaniline	15.563	138	412805	51.009	ng	88
63) Azobenzene	15.827	77	1535895	48.668	ng	94
65) 4,6-Dinitro-2-methylph...	15.627	198	295973	51.314	ng	93
66) n-Nitrosodiphenylamine	15.751	169	1626345	43.771	ng	99
67) 4-Bromophenyl-phenylether	16.433	248	603690	41.877	ng	97
68) Hexachlorobenzene	16.557	284	688802	41.697	ng	98
69) Atrazine	16.710	200	606748	43.943	ng	99
70) Pentachlorophenol	16.904	266	905893	102.952	ng	100
71) Phenanthrene	17.286	178	2946722	45.334	ng	99
72) Anthracene	17.380	178	3027862	46.435	ng	99
73) Carbazole	17.651	167	2770794	49.419	ng	99
74) Di-n-butylphthalate	18.204	149	3120405	46.246	ng	99
75) Fluoranthene	19.257	202	3593230	49.710	ng	97
77) Benzidine	19.433	184	1734373	40.168	ng	98
78) Pyrene	19.610	202	3741284	43.870	ng	98
80) Butylbenzylphthalate	20.486	149	1469602	42.847	ng	99
81) Benzo(a)anthracene	21.321	228	3753370	45.647	ng	98
82) 3,3'-Dichlorobenzidine	21.257	252	1216839	34.255	ng	99
83) Chrysene	21.374	228	3601238	45.180	ng	98
84) Bis(2-ethylhexyl)phtha...	21.251	149	2114202	39.804	ng	97
85) Di-n-octyl phthalate	22.180	149	3624008	40.333	ng	99
87) Indeno(1,2,3-cd)pyrene	26.292	276	5035928	45.408	ng	98
88) Benzo(b)fluoranthene	23.021	252	4111430	45.742	ng	99
89) Benzo(k)fluoranthene	23.068	252	3922550	45.084	ng	99
90) Benzo(a)pyrene	23.651	252	3972536	45.789	ng	99
91) Dibenzo(a,h)anthracene	26.303	278	4247854	45.044	ng	99
92) Benzo(g,h,i)perylene	27.062	276	4125624	44.807	ng	99

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

Quant Time: Jan 19 03:36:59 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 14 18:05:00 2022
Response via : Initial Calibration

