

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
 Data File : BP009572.D
 Acq On : 28 Mar 2022 18:56
 Operator : CG/JU
 Sample : N2140-01MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MHCMSD

Quant Time: Mar 29 03:12:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 06:34:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.764	152	282804	20.000	ng	-0.01
21) Naphthalene-d8	10.557	136	1049025	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	608026	20.000	ng	-0.01
64) Phenanthrene-d10	17.175	188	1174670	20.000	ng	-0.01
76) Chrysene-d12	21.292	240	1104990	20.000	ng	0.00
86) Perylene-d12	23.686	264	1292977	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.370	112	1770904	109.329	ng	0.00
7) Phenol-d6	6.958	99	2197329	100.310	ng	0.00
23) Nitrobenzene-d5	8.922	82	1353729	68.576	ng	-0.01
42) 2,4,6-Tribromophenol	15.910	330	841204	83.840	ng	-0.01
45) 2-Fluorobiphenyl	13.034	172	3052686	69.601	ng	0.00
79) Terphenyl-d14	19.757	244	4304646	69.923	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.305	88	310889	48.457	ng	97
3) Pyridine	3.699	79	854952	49.478	ng	97
4) n-Nitrosodimethylamine	3.605	42	347662	54.632	ng	94
6) Aniline	7.099	93	494492	19.807	ng	99
8) 2-Chlorophenol	7.340	128	896168	48.519	ng	97
9) Benzaldehyde	6.911	77	438325	42.239	ng	98
10) Phenol	6.981	94	1098901	50.004	ng	98
11) bis(2-Chloroethyl)ether	7.193	93	849520	48.844	ng	98
12) 1,3-Dichlorobenzene	7.658	146	1004294	48.175	ng	98
13) 1,4-Dichlorobenzene	7.799	146	1025001	48.090	ng	99
14) 1,2-Dichlorobenzene	8.116	146	975734	47.363	ng	99
15) Benzyl Alcohol	8.016	79	736786	42.186	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.299	45	961832	51.017	ng	97
17) 2-Methylphenol	8.222	107	698701	46.137	ng	98
18) Hexachloroethane	8.840	117	361461	48.390	ng	96
19) n-Nitroso-di-n-propyla...	8.575	70	622314	44.915	ng	99
20) 3+4-Methylphenols	8.552	107	944575	45.361	ng	97
22) Acetophenone	8.587	105	1251319	48.194	ng	99
24) Nitrobenzene	8.963	77	934573	50.116	ng	97
25) Isophorone	9.493	82	1685226	50.048	ng	98
26) 2-Nitrophenol	9.675	139	463419	47.676	ng	96
27) 2,4-Dimethylphenol	9.752	122	722453	52.673	ng	99
28) bis(2-Chloroethoxy)met...	9.975	93	1046730	47.742	ng	99
29) 2,4-Dichlorophenol	10.222	162	800052	47.820	ng	99
30) 1,2,4-Trichlorobenzene	10.428	180	898896	46.374	ng	98
31) Naphthalene	10.610	128	2586770	47.872	ng	100
32) Benzoic acid	9.928	122	485154	45.242	ng	93
33) 4-Chloroaniline	10.734	127	159212	7.221	ng	98
34) Hexachlorobutadiene	10.899	225	579234	44.103	ng	99
35) Caprolactam	11.522	113	227330	40.262	ng	96
36) 4-Chloro-3-methylphenol	11.875	107	780307	43.404	ng	98
37) 2-Methylnaphthalene	12.228	142	1755817	46.312	ng	99
38) 1-Methylnaphthalene	12.446	142	1676371	45.389	ng	100
40) 1,2,4,5-Tetrachloroben...	12.599	216	984895	48.821	ng	99
41) Hexachlorocyclopentadiene	12.581	237	832708	93.475	ng	99
43) 2,4,6-Trichlorophenol	12.846	196	625862	46.806	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
 Data File : BP009572.D
 Acq On : 28 Mar 2022 18:56
 Operator : CG/JU
 Sample : N2140-01MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MHCMSD

Quant Time: Mar 29 03:12:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 06:34:38 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.928	196	673364	46.373	ng	98
46) 1,1'-Biphenyl	13.240	154	2238839	49.358	ng	99
47) 2-Chloronaphthalene	13.281	162	1748527	48.962	ng	99
48) 2-Nitroaniline	13.493	65	456787	46.964	ng	99
49) Acenaphthylene	14.134	152	2810349	50.978	ng	100
50) Dimethylphthalate	13.875	163	2052467	44.549	ng	100
51) 2,6-Dinitrotoluene	13.993	165	452416	46.868	ng	100
52) Acenaphthene	14.475	154	1585890	48.157	ng	99
53) 3-Nitroaniline	14.328	138	209273	20.525	ng	# 96
54) 2,4-Dinitrophenol	14.540	184	496018	82.541	ng	98
55) Dibenzofuran	14.816	168	2524560	45.632	ng	100
56) 4-Nitrophenol	14.657	139	707026	93.028	ng	99
57) 2,4-Dinitrotoluene	14.787	165	600510	45.243	ng	98
58) Fluorene	15.469	166	1982526	45.563	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.051	232	546620	41.524	ng	96
60) Diethylphthalate	15.246	149	1987127	43.095	ng	100
61) 4-Chlorophenyl-phenyle...	15.463	204	1047206	43.690	ng	96
62) 4-Nitroaniline	15.498	138	420150	39.238	ng	98
63) Azobenzene	15.757	77	1880824	46.890	ng	99
65) 4,6-Dinitro-2-methylph...	15.563	198	330278	43.726	ng	93
66) n-Nitrosodiphenylamine	15.681	169	1724214	49.601	ng	98
67) 4-Bromophenyl-phenylether	16.363	248	658830	45.835	ng	95
68) Hexachlorobenzene	16.481	284	745735	45.078	ng	98
69) Atrazine	16.645	200	615708	50.378	ng	99
70) Pentachlorophenol	16.834	266	839765	94.962	ng	99
71) Phenanthrene	17.222	178	3019678	48.023	ng	100
72) Anthracene	17.310	178	3050895	49.142	ng	100
73) Carbazole	17.587	167	2755164	45.378	ng	99
74) Di-n-butylphthalate	18.151	149	3305791	46.463	ng	99
75) Fluoranthene	19.204	202	3503263	46.019	ng	98
77) Benzidine	19.386	184	1066571	39.379	ng	99
78) Pyrene	19.557	202	3590171	49.306	ng	99
80) Butylbenzylphthalate	20.439	149	1419855	45.878	ng	99
81) Benzo(a)anthracene	21.280	228	3504343	48.271	ng	99
82) 3,3'-Dichlorobenzidine	21.210	252	754585	26.906	ng	99
83) Chrysene	21.333	228	3251555	47.300	ng	99
84) Bis(2-ethylhexyl)phtha...	21.216	149	2075238	48.107	ng	100
85) Di-n-octyl phthalate	22.133	149	3540804	47.620	ng	99
87) Indeno(1,2,3-cd)pyrene	26.186	276	4973939	50.705	ng	96
88) Benzo(b)fluoranthene	22.957	252	3789552	49.096	ng	98
89) Benzo(k)fluoranthene	23.004	252	3746346	49.572	ng	98
90) Benzo(a)pyrene	23.580	252	3811617	57.717	ng	98
91) Dibenzo(a,h)anthracene	26.198	278	4332549	52.994	ng	97
92) Benzo(g,h,i)perylene	26.939	276	4356825	54.171	ng	97

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

