

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN072618\
 Data File : BN002175.D
 Acq On : 26 Jul 2018 16:43
 Operator : JU/SJ
 Sample : PB111396BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB111396BS

Quant Time: Jul 27 01:21:02 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN072518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 26 18:27:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	244562	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	890584	20.00	ng	0.00
38) Acenaphthene-d10	14.32	164	436380	20.00	ng	0.00
63) Phenanthrene-d10	17.07	188	775952	20.00	ng	0.00
75) Chrysene-d12	21.27	240	613832	20.00	ng	0.00
86) Perylene-d12	23.49	264	587248	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	2147892	141.39	ng	0.00
7) Phenol-d6	6.88	99	2752029	129.10	ng	0.00
23) Nitrobenzene-d5	8.85	82	1707021	98.30	ng	0.00
41) 2,4,6-Tribromophenol	15.82	330	706444	124.67	ng	0.00
44) 2-Fluorobiphenyl	12.94	172	3344409	99.82	ng	0.00
78) Terphenyl-d14	19.70	244	3218839	109.37	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.29	88	215362	37.763	ng	# 77
3) Pyridine	3.66	79	634732	38.424	ng	# 71
4) n-Nitrosodimethylamine	3.59	42	330948	46.704	ng	# 72
6) Aniline	7.03	93	590269	22.667	ng	98
8) 2-Chlorophenol	7.26	128	751889	42.679	ng	94
9) Benzaldehyde	6.85	77	458907	34.954	ng	95
10) Phenol	6.90	94	925430	42.351	ng	100
11) bis(2-Chloroethyl)ether	7.13	93	714187	40.155	ng	96
12) 1,3-Dichlorobenzene	7.58	146	830247	41.352	ng	99
13) 1,4-Dichlorobenzene	7.73	146	841669	40.978	ng	97
14) 1,2-Dichlorobenzene	8.04	146	802196	40.485	ng	98
15) Benzyl Alcohol	7.93	79	620303	38.561	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.22	45	1172151	40.813	ng	93
17) 2-Methylphenol	8.14	107	606815	40.677	ng	95
18) Hexachloroethane	8.76	117	312469	41.567	ng	89
19) n-Nitroso-di-n-propylamine	8.49	70	560083	36.222	ng	# 98
20) 3+4-Methylphenols	8.46	107	820274	39.410	ng	95
22) Acetophenone	8.51	105	1099672	47.624	ng	97
24) Nitrobenzene	8.89	77	841324	46.746	ng	94
25) Isophorone	9.40	82	1444161	48.432	ng	95
26) 2-Nitrophenol	9.59	139	388127	44.643	ng	97
27) 2,4-Dimethylphenol	9.65	122	637837	52.752	ng	98
28) bis(2-Chloroethoxy)methane	9.89	93	893759	46.402	ng	98
29) 2,4-Dichlorophenol	10.12	162	644921	45.741	ng	95
30) 1,2,4-Trichlorobenzene	10.33	180	731286	45.798	ng	98
31) Naphthalene	10.52	128	2188582	48.693	ng	98
32) Benzoic acid	9.81	122	372030	36.265	ng	91
33) 4-Chloroaniline	10.63	127	128814	6.441	ng	98
34) Hexachlorobutadiene	10.80	225	442294	45.232	ng	97
35) Caprolactam	11.40	113	172476	35.845	ng	# 58
36) 4-Chloro-3-methylphenol	11.76	107	654727	40.838	ng	96
37) 2-Methylnaphthalene	12.13	142	1520216	47.957	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.50	216	738058	50.300	ng	# 98
40) Hexachlorocyclopentadiene	12.48	237	1050554	126.770	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.75	196	464344	47.166	ng	98
43) 2,4,5-Trichlorophenol	12.82	196	483583	46.371	ng	98
45) 1,1'-Biphenyl	13.15	154	1835517	47.800	ng	99
46) 2-Chloronaphthalene	13.19	162	1409689	47.560	ng	99
47) 2-Nitroaniline	13.40	65	413380	43.343	ng	96
48) Acenaphthylene	14.05	152	2226736	49.507	ng	97
49) Dimethylphthalate	13.79	163	1593142	43.950	ng	100
50) 2,6-Dinitrotoluene	13.90	165	345807	42.940	ng	93
51) Acenaphthene	14.39	154	1274093	47.026	ng	99
52) 3-Nitroaniline	14.23	138	103399	11.850	ng	# 88
53) 2,4-Dinitrophenol	14.45	184	356917	83.702	ng	98
54) Dibenzofuran	14.72	168	1958334	44.582	ng	96
55) 4-Nitrophenol	14.56	139	536515	80.090	ng	88
56) 2,4-Dinitrotoluene	14.70	165	455046	41.634	ng	87
57) Fluorene	15.37	166	1529474	46.196	ng	98
58) 2,3,4,6-Tetrachlorophenol	14.95	232	393985	43.929	ng	95
59) Diethylphthalate	15.16	149	1527050	41.389	ng	99
60) 4-Chlorophenyl-phenylether	15.37	204	771421	42.107	ng	96
61) 4-Nitroaniline	15.40	138	315129	35.429	ng	94
62) Azobenzene	15.66	77	1553268	43.875	ng	98
64) 4,6-Dinitro-2-methylphenol	15.46	198	233338	44.541	ng	97
65) n-Nitrosodiphenylamine	15.59	169	1288529	50.043	ng	96
66) 4-Bromophenyl-phenylether	16.27	248	447582	49.273	ng	92
67) Hexachlorobenzene	16.38	284	475616	47.337	ng	95
68) Atrazine	16.55	200	401467	46.311	ng	96
69) Pentachlorophenol	16.73	266	493992	76.380	ng	100
70) Phenanthrene	17.11	178	2135962	50.430	ng	98
71) Anthracene	17.20	178	2154662	51.397	ng	98
72) Carbazole	17.47	167	1847750	43.927	ng	97
73) Di-n-butylphthalate	18.05	149	2258153	46.022	ng	99
74) Fluoranthene	19.13	202	2156759	46.901	ng	98
76) Benzidine	19.32	184	727407	36.334	ng	97
77) Pyrene	19.50	202	2139334	50.538	ng	99
79) Butylbenzylphthalate	20.41	149	884788	48.016	ng	93
80) Benzo(a)anthracene	21.25	228	1884799	50.378	ng	98
81) 3,3'-Dichlorobenzidine	21.19	252	350902	24.075	ng	# 98
82) Chrysene	21.30	228	1738500	48.804	ng	98
83) Bis(2-ethylhexyl)phthalate	21.19	149	1239857	48.251	ng	97
84) Di-n-octyl phthalate	22.07	149	2039263	50.052	ng	98
85) Indeno(1,2,3-cd)pyrene	25.73	276	2203157	63.437	ng	# 95
87) Benzo(b)fluoranthene	22.83	252	1767583	51.001	ng	97
88) Benzo(k)fluoranthene	22.87	252	1721966	50.449	ng	98
89) Benzo(a)pyrene	23.40	252	1700349	52.533	ng	# 97
90) Dibenzo(a,h)anthracene	25.74	278	1836514	60.434	ng	# 97
91) Benzo(g,h,i)perylene	26.40	276	1890589	64.376	ng	# 94

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

