

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082720\
 Data File : BP003134.D
 Acq On : 27 Aug 2020 16:23
 Operator : CG/JU
 Sample : L3806-08MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-14-DMS

Quant Time: Aug 27 17:35:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 15:32:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	223385	20.00	ng	0.00
21) Naphthalene-d8	10.44	136	857683	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	493100	20.00	ng	0.00
64) Phenanthrene-d10	17.06	188	978375	20.00	ng	0.00
76) Chrysene-d12	21.15	240	924445	20.00	ng	0.00
86) Perylene-d12	23.38	264	1110547	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	1337078	105.83	ng	0.00
7) Phenol-d6	6.85	99	1560813	98.55	ng	0.00
23) Nitrobenzene-d5	8.81	82	910389	76.52	ng	0.00
42) 2,4,6-Tribromophenol	15.80	330	683578	102.50	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	2469456	73.34	ng	0.00
79) Terphenyl-d14	19.63	244	2924322	69.61	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.21	88	177892	36.828	ng	99
3) Pyridine	3.59	79	410281	32.275	ng	99
4) n-Nitrosodimethylamine	3.51	42	139613	42.058	ng	92
6) Aniline	6.99	93	543442	31.514	ng	99
8) 2-Chlorophenol	7.23	128	600649	42.643	ng	99
9) Benzaldehyde	6.80	77	261853	35.461	ng	99
10) Phenol	6.88	94	618018	42.105	ng	98
11) bis(2-Chloroethyl)ether	7.09	93	472486	40.670	ng	99
12) 1,3-Dichlorobenzene	7.55	146	676036	39.472	ng	99
13) 1,4-Dichlorobenzene	7.70	146	685657	39.656	ng	99
14) 1,2-Dichlorobenzene	8.01	146	659714	40.246	ng	99
15) Benzyl Alcohol	7.90	79	389157	43.608	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.19	45	373304	37.918	ng	99
17) 2-Methylphenol	8.11	107	443105	42.376	ng	98
18) Hexachloroethane	8.74	117	225389	39.846	ng	96
19) n-Nitroso-di-n-propylamine	8.47	70	286946	40.099	ng	100
20) 3+4-Methylphenols	8.43	107	589426	41.837	ng	98
22) Acetophenone	8.48	105	750485	39.291	ng	99
24) Nitrobenzene	8.85	77	473549	40.525	ng	99
25) Isophorone	9.38	82	878852	41.867	ng	99
26) 2-Nitrophenol	9.56	139	311213	44.683	ng	98
27) 2,4-Dimethylphenol	9.63	122	505036	48.236	ng	97
28) bis(2-Chloroethoxy)methane	9.86	93	590308	37.756	ng	100
29) 2,4-Dichlorophenol	10.09	162	549928	42.335	ng	99
30) 1,2,4-Trichlorobenzene	10.31	180	611593	39.543	ng	99
31) Naphthalene	10.49	128	1762096	39.988	ng	100
32) Benzoic acid	9.78	122	275791	36.664	ng	97
33) 4-Chloroaniline	10.60	127	245074	13.311	ng	99
34) Hexachlorobutadiene	10.79	225	359109	38.687	ng	99
35) Caprolactam	11.36	113	151983	38.934	ng	99
36) 4-Chloro-3-methylphenol	11.74	107	491311	40.508	ng	98
37) 2-Methylnaphthalene	12.11	142	1255755	39.837	ng	99
38) 1-Methylnaphthalene	12.33	142	1181957	39.450	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.49	216	619596	42.394	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082720\
 Data File : BP003134.D
 Acq On : 27 Aug 2020 16:23
 Operator : CG/JU
 Sample : L3806-08MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-14-DMS

Quant Time: Aug 27 17:35:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 15:32:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.48	237	762201	110.324	ng	99
43) 2,4,6-Trichlorophenol	12.73	196	408066	42.463	ng	99
44) 2,4,5-Trichlorophenol	12.80	196	437667	43.056	ng	100
46) 1,1'-Biphenyl	13.13	154	1561214	42.210	ng	100
47) 2-Chloronaphthalene	13.17	162	1196114	39.508	ng	99
48) 2-Nitroaniline	13.37	65	226576	39.970	ng	99
49) Acenaphthylene	14.02	152	1893641	41.364	ng	100
50) Dimethylphthalate	13.77	163	1802003	47.921	ng	100
51) 2,6-Dinitrotoluene	13.87	165	317694	40.651	ng	99
52) Acenaphthene	14.37	154	1110551	39.486	ng	98
53) 3-Nitroaniline	14.20	138	183169	20.687	ng	99
54) 2,4-Dinitrophenol	14.42	184	315731	74.541	ng	97
55) Dibenzofuran	14.70	168	1747595	39.565	ng	99
56) 4-Nitrophenol	14.53	139	548000	86.026	ng	98
57) 2,4-Dinitrotoluene	14.67	165	424164	40.365	ng	98
58) Fluorene	15.36	166	1326418	38.983	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.93	232	379416	41.255	ng	99
60) Diethylphthalate	15.14	149	1392837	37.754	ng	100
61) 4-Chlorophenyl-phenylether	15.36	204	680797	38.745	ng	100
62) 4-Nitroaniline	15.37	138	315169	34.627	ng	98
63) Azobenzene	15.65	77	966988	39.493	ng	99
65) 4,6-Dinitro-2-methylphenol	15.44	198	223801	47.831	ng	99
66) n-Nitrosodiphenylamine	15.57	169	1204272	42.096	ng	99
67) 4-Bromophenyl-phenylether	16.25	248	442822	41.010	ng	99
68) Hexachlorobenzene	16.37	284	514001	39.983	ng	99
69) Atrazine	16.53	200	367141	38.051	ng	100
70) Pentachlorophenol	16.72	266	606833	98.712	ng	99
71) Phenanthrene	17.10	178	2114899	39.893	ng	100
72) Anthracene	17.19	178	2159826	42.568	ng	100
73) Carbazole	17.46	167	1943255	39.907	ng	99
74) Di-n-butylphthalate	18.03	149	2318246	40.499	ng	100
75) Fluoranthene	19.07	202	2410205	39.336	ng	100
77) Benzidine	19.26	184	1230837	58.256	ng	99
78) Pyrene	19.43	202	2456652	40.945	ng	100
80) Butylbenzylphthalate	20.31	149	1011707	40.954	ng	98
81) Benzo(a)anthracene	21.13	228	2321563	40.092	ng	100
82) 3,3'-Dichlorobenzidine	21.07	252	682522	32.043	ng	100
83) Chrysene	21.19	228	2205105	39.773	ng	100
84) Bis(2-ethylhexyl)phthalate	21.08	149	1478234	41.435	ng	100
85) Di-n-octyl phthalate	21.95	149	2579572	42.173	ng	100
87) Indeno(1,2,3-cd)pyrene	25.67	276	3471112	41.912	ng	99
88) Benzo(b)fluoranthene	22.72	252	2552989	36.987	ng	100
89) Benzo(k)fluoranthene	22.76	252	2543995	38.575	ng	99
90) Benzo(a)pyrene	23.29	252	2424832	37.867	ng	99
91) Dibenzo(a,h)anthracene	25.67	278	2881649	42.356	ng	100
92) Benzo(g,h,i)perylene	26.36	276	2997085	43.897	ng	100

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

