

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110719\
 Data File : BG043545.D
 Acq On : 7 Nov 2019 18:51
 Operator : JU
 Sample : K5710-11MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP- 32-DMSD

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:33:02 PM

Quant Time: Nov 08 04:45:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	31701	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	116691	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	82604	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	196047	20.00	ng	0.00
76) Chrysene-d12	21.65	240	212347	20.00	ng	0.00
87) Perylene-d12	24.84	264	245248	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	207000	119.65	ng	0.02
7) Phenol-d6	7.22	99	293644	117.98	ng	0.02
23) Nitrobenzene-d5	9.17	82	174564	77.12	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	172041	109.44	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	438710	73.39	ng	0.00
79) Terphenyl-d14	19.99	244	855848	75.61	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.47	88	23572	34.965	ng	96
3) Pyridine	3.87	79	61681	36.270	ng	97
4) n-Nitrosodimethylamine	3.79	42	36118	42.089	ng	# 96
6) Aniline	7.34	93	53240	18.568	ng	97
8) 2-Chlorophenol	7.59	128	83706	43.832	ng	99
9) Benzaldehyde	7.15	77	42160	34.466	ng	89
10) Phenol	7.24	94	109015	47.930	ng	97
11) bis(2-Chloroethyl)ether	7.43	93	67704	38.501	ng	99
12) 1,3-Dichlorobenzene	7.89	146	93978	39.277	ng	97
13) 1,4-Dichlorobenzene	8.04	146	95847	39.593	ng	97
14) 1,2-Dichlorobenzene	8.36	146	91620	38.762	ng	93
15) Benzyl Alcohol	8.26	79	73930	42.293	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.53	45	93084	36.404	ng	98
17) 2-Methylphenol	8.48	107	69123	41.688	ng	94
18) Hexachloroethane	9.09	117	32986	37.767	ng	98
19) n-Nitroso-di-n-propylamine	8.81	70	60723	37.754	ng	97
20) 3+4-Methylphenols	8.81	107	94070	41.374	ng	# 85
22) Acetophenone	8.83	105	118620	39.694	ng	# 91
24) Nitrobenzene	9.21	77	92505	42.902	ng	# 98
25) Isophorone	9.73	82	169339	40.921	ng	96
26) 2-Nitrophenol	9.93	139	47802	45.921	ng	# 92
27) 2,4-Dimethylphenol	10.00	122	77937	52.237	ng	94
28) bis(2-Chloroethoxy)methane	10.21	93	93001	40.719	ng	98
29) 2,4-Dichlorophenol	10.49	162	87593	45.821	ng	98
30) 1,2,4-Trichlorobenzene	10.67	180	99050	41.110	ng	97
31) Naphthalene	10.86	128	259095	43.743	ng	98
32) Benzoic acid	10.19	122	43436	39.989	ng	96
33) 4-Chloroaniline	10.99	127	28048	11.552	ng	# 92
34) Hexachlorobutadiene	11.15	225	69254	38.883	ng	96
35) Caprolactam	11.75	113	28393m	43.126	ng	
36) 4-Chloro-3-methylphenol	12.13	107	94013	45.086	ng	98
37) 2-Methylnaphthalene	12.47	142	195647	43.049	ng	98
38) 1-Methylnaphthalene	12.68	142	183817	42.509	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.84	216	124567	40.747	ng	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110719\
 Data File : BG043545.D
 Acq On : 7 Nov 2019 18:51
 Operator : JU
 Sample : K5710-11MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP- 32-DMSD

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:33:02 PM

Quant Time: Nov 08 04:45:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.81	237	122287	76.149	nq	98
43) 2,4,6-Trichlorophenol	13.09	196	81538	45.291	nq	98
44) 2,4,5-Trichlorophenol	13.18	196	84423	46.384	nq	95
46) 1,1'-Biphenyl	13.47	154	259213	41.249	nq	100
47) 2-Chloronaphthalene	13.52	162	200461	41.926	nq	97
48) 2-Nitroaniline	13.73	65	60769	42.524	nq	95
49) Acenaphthylene	14.36	152	325377	43.725	nq	99
50) Dimethylphthalate	14.09	163	312049	48.771	nq	99
51) 2,6-Dinitrotoluene	14.22	165	60092	43.231	nq	96
52) Acenaphthene	14.70	154	192243	42.362	nq	97
53) 3-Nitroaniline	14.55	138	21962	16.365	nq	100
54) 2,4-Dinitrophenol	14.76	184	61950	72.157	nq	96
55) Dibenzofuran	15.04	168	318620	43.295	nq	99
56) 4-Nitrophenol	14.90	139	85711	95.305	nq	93
57) 2,4-Dinitrotoluene	15.00	165	85127	42.853	nq	# 98
58) Fluorene	15.68	166	268537	43.507	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.27	232	85740	44.338	nq	96
60) Diethylphthalate	15.45	149	263881	41.046	nq	98
61) 4-Chlorophenyl-phenylether	15.67	204	149959	41.646	nq	96
62) 4-Nitroaniline	15.72	138	55988	40.396	nq	98
63) Azobenzene	15.97	77	223630	42.981	nq	95
65) 4,6-Dinitro-2-methylphenol	15.77	198	51947	45.025	nq	94
66) n-Nitrosodiphenylamine	15.89	169	236707	42.766	nq	98
67) 4-Bromophenyl-phenylether	16.57	248	106665	40.429	nq	99
68) Hexachlorobenzene	16.70	284	120311	39.726	nq	99
69) Atrazine	16.84	200	102342	53.213	nq	96
70) Pentachlorophenol	17.05	266	108087	78.325	nq	94
71) Phenanthrene	17.43	178	426463	42.480	nq	100
72) Anthracene	17.52	178	443671	44.172	nq	99
73) Carbazole	17.79	167	399999	43.976	nq	98
74) Di-n-butylphthalate	18.34	149	466900	42.305	nq	99
75) Fluoranthene	19.43	202	572731	43.760	nq	99
77) Benzidine	19.62	184	189418	44.323	nq	99
78) Pyrene	19.80	202	576306	43.845	nq	99
80) Butylbenzylphthalate	20.68	149	216597	43.496	nq	96
81) Benzo(a)anthracene	21.63	228	593654	44.631	nq	98
82) 3,3'-Dichlorobenzidine	21.55	252	105677	20.541	nq	99
83) Chrysene	21.70	228	565814	42.813	nq	98
84) Bis(2-ethylhexyl)phthalate	21.53	149	312423	44.166	nq	99
85) Di-n-octyl phthalate	22.73	149	521592	43.462	nq	100
86) Indeno(1,2,3-cd)pyrene	28.49	276	715677	43.141	nq	99
88) Benzo(b)fluoranthene	23.83	252	595562	42.877	nq	99
89) Benzo(k)fluoranthene	23.90	252	615529	44.019	nq	98
90) Benzo(a)pyrene	24.70	252	562380	42.399	nq	97
91) Dibenzo(a,h)anthracene	28.54	278	614179	45.270	nq	99
92) Benzo(g,h,i)perylene	29.62	276	605903	45.275	nq	99

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

