

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081017\
 Data File : BF097572.D
 Acq On : 10 Aug 2017 21:51
 Operator : SJ/JU
 Sample : I4689-01MSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRAP-ROCK-FINESMSD

Manual Integrations
 APPROVED

sohil
 8/14/2017 6:34:21 PM

Quant Time: Aug 11 11:59:05 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.39	152	107152	20.00	ng	-0.02
21) Naphthalene-d8	7.67	136	453299	20.00	ng	-0.02
38) Acenaphthene-d10	9.42	164	188997	20.00	ng	-0.02
63) Phenanthrene-d10	10.89	188	299230	20.00	ng	-0.02
75) Chrysene-d12	13.52	240	153406	20.00	ng	-0.02
86) Perylene-d12	14.86	264	188563	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	4.99	112	604493	85.04	ng	-0.01
7) Phenol-d6	6.07	99	730449	90.02	ng	-0.01
23) Nitrobenzene-d5	6.97	82	451058	58.37	ng	-0.02
41) 2,4,6-Tribromophenol	10.21	330	123947	83.00	ng	-0.02
44) 2-Fluorobiphenyl	8.75	172	671389	60.29	ng	-0.02
78) Terphenyl-d14	12.47	244	432202	57.44	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.89	88	74269	25.038	ng	# 73
3) Pyridine	2.50	79	206296	22.713	ng	# 61
4) n-Nitrosodimethylamine	2.45	42	129814	25.798	ng	80
6) Aniline	6.06	93	259388	25.402	ng	96
8) 2-Chlorophenol	6.19	128	226114	29.750	ng	94
9) Benzaldehyde	5.95	77	94445	19.663	ng	94
10) Phenol	6.09	94	321863	33.440	ng	98
11) bis(2-Chloroethyl)ether	6.14	93	225392	29.608	ng	91
12) 1,3-Dichlorobenzene	6.33	146	220972	26.978	ng	98
13) 1,4-Dichlorobenzene	6.41	146	231593	28.531	ng	95
14) 1,2-Dichlorobenzene	6.56	146	204907	28.911	ng	99
15) Benzyl Alcohol	6.56	79	187711	36.537	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.68	45	470138	30.791	ng	90
17) 2-Methylphenol	6.69	107	182835	31.169	ng	# 90
18) Hexachloroethane	6.89	117	79777	26.864	ng	# 77
19) n-Nitroso-di-n-propylamine	6.83	70	161388	30.215	ng	# 79
20) 3+4-Methylphenols	6.86	107	243971	31.845	ng	# 64
22) Acetophenone	6.82	105	309609	28.441	ng	# 99
24) Nitrobenzene	6.99	77	236749	29.419	ng	96
25) Isophorone	7.23	82	411759	27.210	ng	97
26) 2-Nitrophenol	7.30	139	121962	28.012	ng	# 81
27) 2,4-Dimethylphenol	7.37	122	196070	27.300	ng	97
28) bis(2-Chloroethoxy)methane	7.45	93	282185	28.575	ng	98
29) 2,4-Dichlorophenol	7.57	162	185352	29.957	ng	97
30) 1,2,4-Trichlorobenzene	7.62	180	159839	27.103	ng	98
31) Naphthalene	7.70	128	614466	28.756	ng	100
33) 4-Chloroaniline	7.77	127	191164	21.987	ng	99
34) Hexachlorobutadiene	7.82	225	81330	26.013	ng	98
35) Caprolactam	8.14	113	47977	24.182	ng	# 51
36) 4-Chloro-3-methylphenol	8.27	107	193732	28.701	ng	# 84
37) 2-Methylnaphthalene	8.39	142	407129	30.093	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.56	216	146123	28.976	ng	99
40) Hexachlorocyclopentadiene	8.54	237	19892	16.942	ng	94
42) 2,4,6-Trichlorophenol	8.69	196	95639	26.170	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.74	196	91982	25.147	nq	98
45) 1,1'-Biphenyl	8.85	154	454654	28.164	nq	98
46) 2-Chloronaphthalene	8.87	162	345941	29.121	nq #	94
47) 2-Nitroaniline	8.99	65	140280	32.538	nq	92
48) Acenaphthylene	9.28	152	541663	29.706	nq	99
49) Dimethylphthalate	9.16	163	546947	38.109	nq	99
50) 2,6-Dinitrotoluene	9.23	165	88218	28.981	nq #	81
51) Acenaphthene	9.45	154	315491	29.374	nq	99
52) 3-Nitroaniline	9.40	138	95619	25.307	nq #	93
54) Dibenzofuran	9.63	168	478106	33.420	nq	95
55) 4-Nitrophenol	9.62	139	34202m	15.222	nq	
56) 2,4-Dinitrotoluene	9.65	165	120000	34.292	nq #	80
57) Fluorene	9.96	166	342449	31.849	nq	100
58) 2,3,4,6-Tetrachlorophenol	9.76	232	67609	26.770	nq	96
59) Diethylphthalate	9.86	149	405383	30.788	nq	98
60) 4-Chlorophenyl-phenylether	9.96	204	141880	31.954	nq	95
61) 4-Nitroaniline	10.02	138	90793	25.290	nq	93
62) Azobenzene	10.12	77	405055	33.160	nq	92
64) 4,6-Dinitro-2-methylphenol	10.07	198	10115	5.440	nq #	65
65) n-Nitrosodiphenylamine	10.09	169	321613	30.890	nq	98
66) 4-Bromophenyl-phenylether	10.45	248	89600	30.004	nq	93
67) Hexachlorobenzene	10.52	284	91960	27.461	nq #	90
68) Atrazine	10.62	200	93530	31.328	nq	93
69) Pentachlorophenol	10.73	266	28602	20.643	nq	96
70) Phenanthrene	10.92	178	475233	29.641	nq	99
71) Anthracene	10.97	178	494353	30.105	nq	99
72) Carbazole	11.14	167	449154	27.812	nq	98
73) Di-n-butylphthalate	11.47	149	599724	30.042	nq	98
74) Fluoranthene	12.10	202	398091	25.345	nq	98
76) Benzidine	12.23	184	203089	35.679	nq #	93
77) Pyrene	12.32	202	385108	30.664	nq	99
79) Butylbenzylphthalate	12.96	149	189362	29.642	nq #	83
80) Benzo(a)anthracene	13.51	228	261678	26.363	nq	100
81) 3,3'-Dichlorobenzidine	13.48	252	95770	25.585	nq #	96
82) Chrysene	13.55	228	262065	27.038	nq	100
83) Bis(2-ethylhexyl)phthalate	13.52	149	216916	26.006	nq	99
84) Di-n-octyl phthalate	14.13	149	404275	26.411	nq #	93
85) Indeno(1,2,3-cd)pyrene	16.04	276	238910	28.746	nq	99
87) Benzo(b)fluoranthene	14.52	252	315738	27.855	nq	98
88) Benzo(k)fluoranthene	14.54	252	257985m	23.885	nq	
89) Benzo(a)pyrene	14.81	252	301163	29.031	nq	98
90) Dibenzo(a,h)anthracene	16.04	278	203487	23.919	nq	98
91) Benzo(g,h,i)perylene	16.39	276	195225	21.883	nq	97

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

