

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102417\
 Data File : BF099799.D
 Acq On : 24 Oct 2017 20:24
 Operator : SJ/JU
 Sample : I5956-01MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-M-6(50-52)MSD

Manual Integrations
 APPROVED

Sohil
 10/25/2017 6:25:52 PM

Quant Time: Oct 25 08:32:47 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 23:39:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	89305	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	363079	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	152333	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	221461	20.00	ng	0.00
75) Chrysene-d12	13.99	240	156426	20.00	ng	0.00
86) Perylene-d12	15.45	264	160435	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	610554	107.33	ng	0.01
7) Phenol-d6	6.46	99	716900	106.29	ng	0.00
23) Nitrobenzene-d5	7.37	82	412793	73.20	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	126240	90.94	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	769218	77.91	ng	0.00
78) Terphenyl-d14	12.92	244	468832	65.50	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.58	88	79263	31.554	ng	# 90
3) Pyridine	3.36	79	192099	31.801	ng	# 89
4) n-Nitrosodimethylamine	3.30	42	80201	37.164	ng	# 75
6) Aniline	6.47	93	255775	29.247	ng	# 91
8) 2-Chlorophenol	6.60	128	236514	39.012	ng	91
9) Benzaldehyde	6.36	77	134452	33.359	ng	91
10) Phenol	6.47	94	289296	39.066	ng	86
11) bis(2-Chloroethyl)ether	6.55	93	218489	38.207	ng	94
12) 1,3-Dichlorobenzene	6.75	146	250539	36.805	ng	96
13) 1,4-Dichlorobenzene	6.82	146	255906	37.041	ng	98
14) 1,2-Dichlorobenzene	6.97	146	231803	36.815	ng	97
15) Benzyl Alcohol	6.96	79	162928	37.984	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.08	45	273272	43.018	ng	72
17) 2-Methylphenol	7.07	107	187273	36.929	ng	97
18) Hexachloroethane	7.31	117	79568	34.521	ng	98
19) n-Nitroso-di-n-propylamine	7.22	70	143928	36.414	ng	90
20) 3+4-Methylphenols	7.22	107	233490	39.543	ng	# 68
22) Acetophenone	7.22	105	299634	36.245	ng	# 98
24) Nitrobenzene	7.40	77	220473	38.799	ng	# 88
25) Isophorone	7.63	82	413485	40.405	ng	97
26) 2-Nitrophenol	7.71	139	117849	36.210	ng	# 86
27) 2,4-Dimethylphenol	7.75	122	203698	42.478	ng	97
28) bis(2-Chloroethoxy)methane	7.84	93	274075	38.242	ng	99
29) 2,4-Dichlorophenol	7.96	162	194825	39.109	ng	95
30) 1,2,4-Trichlorobenzene	8.03	180	199229	37.431	ng	97
31) Naphthalene	8.11	128	649384	37.052	ng	100
33) 4-Chloroaniline	8.17	127	182897	25.272	ng	# 94
34) Hexachlorobutadiene	8.22	225	97254	35.523	ng	98
35) Caprolactam	8.54	113	58764m	37.511	ng	
36) 4-Chloro-3-methylphenol	8.66	107	188533	37.080	ng	92
37) 2-Methylnaphthalene	8.80	142	443369	37.750	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	179566	36.497	ng	# 98
40) Hexachlorocyclopentadiene	8.95	237	7084	18.259	ng	99
42) 2,4,6-Trichlorophenol	9.09	196	120192	40.387	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.13	196	121714	38.540	nq	95
45) 1,1'-Biphenyl	9.26	154	506814	36.777	nq	98
46) 2-Chloronaphthalene	9.29	162	394130	39.381	nq	96
47) 2-Nitroaniline	9.40	65	104353	39.728	nq	88
48) Acenaphthylene	9.71	152	561730	36.442	nq	99
49) Dimethylphthalate	9.57	163	481975	39.953	nq	99
50) 2,6-Dinitrotoluene	9.64	165	96112	37.899	nq #	85
51) Acenaphthene	9.88	154	334148	35.478	nq	99
52) 3-Nitroaniline	9.82	138	94351	31.575	nq	86
54) Dibenzofuran	10.05	168	484165	36.417	nq	98
55) 4-Nitrophenol	9.99	139	99089	63.465	nq #	77
56) 2,4-Dinitrotoluene	10.05	165	108568	35.548	nq #	96
57) Fluorene	10.40	166	383451	34.834	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.17	232	78192	35.313	nq	93
59) Diethylphthalate	10.26	149	432477	36.711	nq	98
60) 4-Chlorophenyl-phenylether	10.38	204	182608	35.248	nq	96
61) 4-Nitroaniline	10.43	138	90841	31.864	nq	82
62) Azobenzene	10.55	77	361799	36.637	nq	89
64) 4,6-Dinitro-2-methylphenol	10.47	198	5115	3.674	nq	82
65) n-Nitrosodiphenylamine	10.50	169	344984	42.199	nq	98
66) 4-Bromophenyl-phenylether	10.87	248	98846	42.397	nq #	89
67) Hexachlorobenzene	10.94	284	92308	38.700	nq	94
68) Atrazine	11.03	200	100031	40.734	nq	98
69) Pentachlorophenol	11.14	266	61267	60.113	nq	99
70) Phenanthrene	11.36	178	459484	36.764	nq	98
71) Anthracene	11.41	178	477389	37.630	nq	98
72) Carbazole	11.57	167	423559	35.504	nq	98
73) Di-n-butylphthalate	11.89	149	593124	38.979	nq	99
74) Fluoranthene	12.55	202	413531	31.838	nq	98
76) Benzidine	12.68	184	323359	47.242	nq	98
77) Pyrene	12.78	202	416803	35.042	nq	99
79) Butylbenzylphthalate	13.39	149	198740	33.931	nq	91
80) Benzo(a)anthracene	13.97	228	364078	36.715	nq	98
81) 3,3'-Dichlorobenzidine	13.93	252	131898	36.643	nq	99
82) Chrysene	14.01	228	350878	37.637	nq	99
83) Bis(2-ethylhexyl)phthalate	13.94	149	281303	34.199	nq #	99
84) Di-n-octyl phthalate	14.55	149	503632	35.674	nq #	93
85) Indeno(1,2,3-cd)pyrene	16.93	276	305820	35.733	nq	96
87) Benzo(b)fluoranthene	15.02	252	379410	37.451	nq	96
88) Benzo(k)fluoranthene	15.05	252	346744	36.297	nq	99
89) Benzo(a)pyrene	15.39	252	346486	38.075	nq	98
90) Dibenzo(a,h)anthracene	16.93	278	265934	32.888	nq	96
91) Benzo(g,h,i)perylene	17.37	276	241126	30.479	nq	98

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

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