

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
 Data File : BF104704.D
 Acq On : 23 Apr 2018 12:54
 Operator : JU/SJ
 Sample : J2461-03MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 N-P001-S002-0001-01MSD

Quant Time: Apr 23 16:03:43 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 23 15:38:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	109250	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	407281	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	163116	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	277098	20.00	ng	0.00
75) Chrysene-d12	13.99	240	179772	20.00	ng	0.00
86) Perylene-d12	15.46	264	128680	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	656848	91.93	ng	0.02
7) Phenol-d6	6.46	99	755826	86.10	ng	0.02
23) Nitrobenzene-d5	7.37	82	460887	73.89	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	156818	92.68	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	784712	76.00	ng	0.00
78) Terphenyl-d14	12.93	244	701269	88.93	ng	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.52	88	98688	29.643	ng	89
3) Pyridine	3.29	79	261399	27.926	ng	88
4) n-Nitrosodimethylamine	3.23	42	102847	31.432	ng	# 82
6) Aniline	6.47	93	49873	4.089	ng	# 1
8) 2-Chlorophenol	6.60	128	251127	33.300	ng	92
9) Benzaldehyde	6.36	77	146766	30.565	ng	100
10) Phenol	6.47	94	300304	32.240	ng	98
11) bis(2-Chloroethyl)ether	6.55	93	256891	34.560	ng	94
12) 1,3-Dichlorobenzene	6.75	146	269899	31.592	ng	99
13) 1,4-Dichlorobenzene	6.83	146	274593	32.243	ng	98
14) 1,2-Dichlorobenzene	6.98	146	253940	32.177	ng	99
15) Benzyl Alcohol	6.96	79	190604	28.586	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.09	45	279523	28.002	ng	61
17) 2-Methylphenol	7.07	107	190312	29.032	ng	# 90
18) Hexachloroethane	7.32	117	61710	20.323	ng	97
19) n-Nitroso-di-n-propylamine	7.22	70	163143	28.439	ng	95
20) 3+4-Methylphenols	7.23	107	246500	30.320	ng	96
22) Acetophenone	7.22	105	337463	33.655	ng	96
24) Nitrobenzene	7.39	77	239680	34.805	ng	99
25) Isophorone	7.63	82	428197	32.465	ng	99
26) 2-Nitrophenol	7.71	139	97352	34.726	ng	93
27) 2,4-Dimethylphenol	7.75	122	202971	34.869	ng	99
28) bis(2-Chloroethoxy)methane	7.84	93	286778	35.562	ng	99
29) 2,4-Dichlorophenol	7.96	162	182685	32.892	ng	97
30) 1,2,4-Trichlorobenzene	8.03	180	198177	32.513	ng	99
31) Naphthalene	8.12	128	662435	32.664	ng	99
33) 4-Chloroaniline	8.17	127	47254	5.439	ng	98
34) Hexachlorobutadiene	8.23	225	111104	33.278	ng	99
35) Caprolactam	8.53	113	44880	23.163	ng	# 82
36) 4-Chloro-3-methylphenol	8.66	107	178195	29.921	ng	100
37) 2-Methylnaphthalene	8.80	142	418207	31.880	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	175894	37.670	ng	98
40) Hexachlorocyclopentadiene	8.95	237	5782	2.212	ng	93
42) 2,4,6-Trichlorophenol	9.09	196	112083	34.579	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.14	196	115990	31.978	ng	96
45) 1,1'-Biphenyl	9.27	154	480519	35.067	ng	99
46) 2-Chloronaphthalene	9.30	162	372698	34.434	ng	97
47) 2-Nitroaniline	9.40	65	116136	43.388	ng	97
48) Acenaphthylene	9.71	152	511636	30.129	ng	99
49) Dimethylphthalate	9.57	163	478070	37.166	ng	99
50) 2,6-Dinitrotoluene	9.63	165	68183	28.647	ng	97
51) Acenaphthene	9.88	154	316657	30.562	ng	99
52) 3-Nitroaniline	9.80	138	45753	16.420	ng	96
53) 2,4-Dinitrophenol	9.93	184	719	6.229	ng	# 30
54) Dibenzofuran	10.06	168	473168	32.769	ng	99
55) 4-Nitrophenol	9.99	139	119023	54.377	ng	95
56) 2,4-Dinitrotoluene	10.05	165	73317	23.483	ng	# 93
57) Fluorene	10.40	166	362590	34.499	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.18	232	89269	32.989	ng	93
59) Diethylphthalate	10.27	149	404180	31.550	ng	98
60) 4-Chlorophenyl-phenylether	10.39	204	169477	36.208	ng	99
61) 4-Nitroaniline	10.42	138	54697	20.076	ng	97
62) Azobenzene	10.55	77	352430	28.795	ng	93
64) 4,6-Dinitro-2-methylphenol	10.45	198	1644	8.379	ng	# 61
65) n-Nitrosodiphenylamine	10.51	169	312293	32.504	ng	99
66) 4-Bromophenyl-phenylether	10.88	248	98804	33.522	ng	91
67) Hexachlorobenzene	10.95	284	110752	33.394	ng	# 88
68) Atrazine	11.04	200	92647	31.947	ng	99
69) Pentachlorophenol	11.15	266	90261	43.992	ng	94
70) Phenanthrene	11.36	178	516984	33.502	ng	98
71) Anthracene	11.42	178	520851	33.204	ng	99
72) Carbazole	11.57	167	465257	30.353	ng	100
73) Di-n-butylphthalate	11.90	149	670098	35.788	ng	98
74) Fluoranthene	12.56	202	526641	32.664	ng	98
76) Benzidine	12.67	184	1401	Below	Cal	# 1
77) Pyrene	12.79	202	520382	39.052	ng	99
79) Butylbenzylphthalate	13.40	149	315947	50.328	ng	97
80) Benzo(a)anthracene	13.98	228	365707	34.052	ng	100
82) Chrysene	14.02	228	347634	31.513	ng	100
83) Bis(2-ethylhexyl)phthalate	13.96	149	508631	62.990	ng	99
84) Di-n-octyl phthalate	14.58	149	523151	38.774	ng	97
85) Indeno(1,2,3-cd)pyrene	16.94	276	280346	29.916	ng	96
87) Benzo(b)fluoranthene	15.03	252	282117	34.713	ng	# 97
88) Benzo(k)fluoranthene	15.06	252	259221	33.784	ng	98
89) Benzo(a)pyrene	15.40	252	244005	33.555	ng	96
90) Dibenzo(a,h)anthracene	16.95	278	234394	39.246	ng	99
91) Benzo(g,h,i)perylene	17.39	276	233879	40.420	ng	94

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

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