

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080318\
 Data File : BF107941.D
 Acq On : 3 Aug 2018 16:50
 Operator : JU/SJ
 Sample : J4285-05MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KG-PIT-2MSD

Manual Integrations
 APPROVED

Sohil
 8/6/2018 5:05:28 PM

Quant Time: Aug 03 17:52:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	127569	20.00	ng	-0.01
21) Naphthalene-d8	8.24	136	567336	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	246578	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	463514	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	309250	20.00	ng	-0.01
86) Perylene-d12	15.68	264	366059	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	552297	73.58	ng	0.00
7) Phenol-d6	6.60	99	751622	75.98	ng	0.00
23) Nitrobenzene-d5	7.52	82	511883	51.98	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	233600	69.72	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	1041507	58.86	ng	-0.01
78) Terphenyl-d14	13.07	244	848083	59.47	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.84	88	72318	23.850	ng	# 83
3) Pyridine	3.68	79	154848m	21.538	ng	
4) n-Nitrosodimethylamine	3.65	42	64808m	16.969	ng	
6) Aniline	6.65	93	181049	17.596	ng	# 68
8) 2-Chlorophenol	6.75	128	252125	26.741	ng	94
9) Benzaldehyde	6.61	77	110717	18.550	ng	98
10) Phenol	6.62	94	339150	28.748	ng	90
11) bis(2-Chloroethyl)ether	6.68	93	296884	28.301	ng	96
12) 1,3-Dichlorobenzene	6.90	146	251024	25.384	ng	96
13) 1,4-Dichlorobenzene	6.97	146	292459	26.568	ng	99
14) 1,2-Dichlorobenzene	7.12	146	278691	27.273	ng	99
15) Benzyl Alcohol	7.12	79	142617	23.195	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.22	45	409466	27.149	ng	66
17) 2-Methylphenol	7.22	107	173954	24.993	ng	# 65
18) Hexachloroethane	7.45	117	99859	25.238	ng	94
19) n-Nitroso-di-n-propylamine	7.35	70	177685	28.315	ng	96
20) 3+4-Methylphenols	7.38	107	202340	23.678	ng	# 72
22) Acetophenone	7.37	105	366186	26.828	ng	96
24) Nitrobenzene	7.54	77	226434	21.928	ng	98
25) Isophorone	7.77	82	498634	30.846	ng	99
26) 2-Nitrophenol	7.86	139	135932	26.230	ng	97
27) 2,4-Dimethylphenol	7.90	122	216519	31.206	ng	96
28) bis(2-Chloroethoxy)methane	7.98	93	290164	27.237	ng	96
29) 2,4-Dichlorophenol	8.11	162	201894	25.725	ng	98
30) 1,2,4-Trichlorobenzene	8.17	180	233314	26.059	ng	98
31) Naphthalene	8.25	128	778658	29.526	ng	99
32) Benzoic acid	7.90	122	216519	46.171	ng	# 15
33) 4-Chloroaniline	8.34	127	187063	16.606	ng	97
34) Hexachlorobutadiene	8.36	225	137324	24.820	ng	97
35) Caprolactam	8.67	113	50172	21.236	ng	# 86
36) 4-Chloro-3-methylphenol	8.80	107	208667	24.481	ng	97
37) 2-Methylnaphthalene	8.95	142	498550	30.095	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.11	216	245057	29.149	ng	# 96
40) Hexachlorocyclopentadiene	9.09	237	94710	31.160	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.24	196	123282	26.525	nq	99
43) 2,4,5-Trichlorophenol	9.29	196	158846	28.390	nq	96
45) 1,1'-Biphenyl	9.41	154	652904	29.096	nq	97
46) 2-Chloronaphthalene	9.44	162	479312	28.325	nq	97
47) 2-Nitroaniline	9.56	65	144688	26.286	nq	95
48) Acenaphthylene	9.85	152	766370	29.213	nq	100
49) Dimethylphthalate	9.71	163	647409	34.961	nq	99
50) 2,6-Dinitrotoluene	9.78	165	113061	28.094	nq	90
51) Acenaphthene	10.02	154	451888	29.653	nq	98
52) 3-Nitroaniline	9.98	138	104861	20.678	nq	96
53) 2,4-Dinitrophenol	10.13	184	9510	12.451	nq #	68
54) Dibenzofuran	10.20	168	639571	27.153	nq	97
55) 4-Nitrophenol	10.20	139	367067	145.207	nq #	20
56) 2,4-Dinitrotoluene	10.20	165	140144	26.495	nq #	75
57) Fluorene	10.54	166	519924	28.403	nq	95
58) 2,3,4,6-Tetrachlorophenol	10.32	232	123831	24.096	nq #	77
59) Diethylphthalate	10.40	149	626579	31.470	nq	99
60) 4-Chlorophenyl-phenylether	10.53	204	257452	25.817	nq	92
61) 4-Nitroaniline	10.61	138	101220	22.492	nq	99
62) Azobenzene	10.69	77	477426	25.747	nq	95
64) 4,6-Dinitro-2-methylphenol	10.61	198	23913	9.483	nq #	77
65) n-Nitrosodiphenylamine	10.65	169	447791	29.340	nq	99
66) 4-Bromophenyl-phenylether	11.02	248	156365	28.943	nq #	89
67) Hexachlorobenzene	11.09	284	163738	27.331	nq #	89
68) Atrazine	11.18	200	121480	26.893	nq	95
69) Pentachlorophenol	11.30	266	77046	26.125	nq	95
70) Phenanthrene	11.51	178	627896	28.529	nq	98
71) Anthracene	11.57	178	774722	30.518	nq	99
72) Carbazole	11.74	167	641589	25.575	nq	99
73) Di-n-butylphthalate	12.03	149	766255	27.997	nq	100
74) Fluoranthene	12.71	202	663528	25.662	nq	96
76) Benzidine	12.85	184	236325	23.635	nq	97
77) Pyrene	12.94	202	620795	27.878	nq	97
79) Butylbenzylphthalate	13.54	149	251405	25.781	nq	93
80) Benzo(a)anthracene	14.13	228	480081	27.405	nq	99
81) 3,3'-Dichlorobenzidine	14.10	252	169139	24.907	nq #	97
82) Chrysene	14.17	228	562168	29.753	nq	98
83) Bis(2-ethylhexyl)phthalate	14.09	149	329771	26.900	nq	99
84) Di-n-octyl phthalate	14.71	149	587804	29.872	nq	97
85) Indeno(1,2,3-cd)pyrene	17.26	276	647026	45.020	nq	93
87) Benzo(b)fluoranthene	15.22	252	456723	25.298	nq	98
88) Benzo(k)fluoranthene	15.25	252	646025	29.145	nq	97
89) Benzo(a)pyrene	15.61	252	556125	29.838	nq	98
90) Dibenzo(a,h)anthracene	17.27	278	542930	33.395	nq #	95
91) Benzo(g,h,i)perylene	17.74	276	507657	32.271	nq #	93

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

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