

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022318\
 Data File : BF103193.D
 Acq On : 23 Feb 2018 12:25
 Operator : SJ/JU
 Sample : J1161-03
 Misc : MDL-S-8ppm
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MDL-SOIL-03-QT1-2018

Manual Integrations
 APPROVED

Sohil
 2/26/2018 9:54:36 AM

Quant Time: Feb 23 12:55:09 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	173095	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	763415	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	353890	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	604884	20.00	ng	0.00
75) Chrysene-d12	14.07	240	453170	20.00	ng	0.02
86) Perylene-d12	15.60	264	433284	20.00	ng	0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	1282471	112.06	ng	0.00
7) Phenol-d6	6.50	99	1569851	112.10	ng	0.00
23) Nitrobenzene-d5	7.44	82	1177826	88.11	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	343717	114.75	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1753046	85.42	ng	0.00
78) Terphenyl-d14	12.99	244	1406746	74.62	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.65	88	45502	7.528	ng	94
3) Pyridine	3.53	79	120229m	7.450	ng	
4) n-Nitrosodimethylamine	3.38	42	49337m	7.581	ng	
6) Aniline	6.54	93	169147	8.369	ng	98
8) 2-Chlorophenol	6.66	128	96088	8.082	ng	97
9) Benzaldehyde	6.43	77	56939	4.914	ng	100
10) Phenol	6.52	94	131503	8.089	ng	96
11) bis(2-Chloroethyl)ether	6.60	93	104454	8.465	ng	91
12) 1,3-Dichlorobenzene	6.81	146	113212	8.332	ng	99
13) 1,4-Dichlorobenzene	6.89	146	114761	8.425	ng	99
14) 1,2-Dichlorobenzene	7.04	146	104016	8.281	ng	96
15) Benzyl Alcohol	7.01	79	101899	9.656	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.14	45	176047	8.308	ng	99
17) 2-Methylphenol	7.12	107	89223	8.258	ng	98
18) Hexachloroethane	7.38	117	41423	8.353	ng	97
19) n-Nitroso-di-n-propylamine	7.27	70	81277	9.325	ng	92
20) 3+4-Methylphenols	7.27	107	117038	8.886	ng	# 37
22) Acetophenone	7.27	105	158283	8.883	ng	# 90
24) Nitrobenzene	7.45	77	121552	8.259	ng	95
25) Isophorone	7.69	82	211620	8.038	ng	97
26) 2-Nitrophenol	7.77	139	52080	7.517	ng	96
27) 2,4-Dimethylphenol	7.80	122	78250	7.389	ng	98
28) bis(2-Chloroethoxy)methane	7.89	93	126981	8.203	ng	98
29) 2,4-Dichlorophenol	8.01	162	78934	7.785	ng	97
30) 1,2,4-Trichlorobenzene	8.09	180	86920	8.050	ng	99
31) Naphthalene	8.17	128	312720	8.367	ng	99
32) Benzoic acid	7.89	122	45954	11.920	ng	98
33) 4-Chloroaniline	8.23	127	128468	7.965	ng	98
34) Hexachlorobutadiene	8.29	225	45575	8.105	ng	98
35) Caprolactam	8.57	113	26493	7.434	ng	# 85
36) 4-Chloro-3-methylphenol	8.70	107	90579	7.920	ng	94
37) 2-Methylnaphthalene	8.86	142	208891	8.648	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	79696	8.238	ng	99
40) Hexachlorocyclopentadiene	9.02	237	7918	10.946	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	50484	7.755	ng	98
43) 2,4,5-Trichlorophenol	9.19	196	55137	7.364	ng	99
45) 1,1'-Biphenyl	9.33	154	261751	8.827	ng	98
46) 2-Chloronaphthalene	9.36	162	194876	8.307	ng	97
47) 2-Nitroaniline	9.45	65	67876	7.811	ng	88
48) Acenaphthylene	9.77	152	310626	8.605	ng	98
49) Dimethylphthalate	9.62	163	233220	8.215	ng	99
50) 2,6-Dinitrotoluene	9.69	165	47881	8.037	ng	96
51) Acenaphthene	9.94	154	181873	8.641	ng	99
52) 3-Nitroaniline	9.86	138	58453	7.733	ng	99
53) 2,4-Dinitrophenol	9.98	184	8910	13.264	ng #	12
54) Dibenzofuran	10.12	168	262585	9.088	ng	99
55) 4-Nitrophenol	10.04	139	27337	5.835	ng	95
56) 2,4-Dinitrotoluene	10.10	165	61225	8.126	ng	95
57) Fluorene	10.46	166	209512	8.512	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.23	232	38942	7.742	ng	93
59) Diethylphthalate	10.32	149	232378	8.558	ng	99
60) 4-Chlorophenyl-phenylether	10.45	204	91739	8.789	ng	97
61) 4-Nitroaniline	10.47	138	55250	7.318	ng	95
62) Azobenzene	10.61	77	235415	8.518	ng	98
64) 4,6-Dinitro-2-methylphenol	10.51	198	21163	10.103	ng	85
65) n-Nitrosodiphenylamine	10.56	169	178730	8.486	ng	99
66) 4-Bromophenyl-phenylether	10.94	248	50905	8.079	ng	95
67) Hexachlorobenzene	11.01	284	55598	8.129	ng	92
68) Atrazine	11.09	200	48906	7.726	ng	98
69) Pentachlorophenol	11.20	266	19495	5.895	ng	93
70) Phenanthrene	11.42	178	287774	8.645	ng	98
71) Anthracene	11.47	178	293105	8.587	ng	99
72) Carbazole	11.63	167	285171	8.389	ng	100
73) Di-n-butylphthalate	11.95	149	357981	8.416	ng	99
74) Fluoranthene	12.62	202	278605	8.329	ng	99
76) Benzidine	12.73	184	72934	4.305	ng #	93
77) Pyrene	12.84	202	289383	8.190	ng	99
79) Butylbenzylphthalate	13.46	149	142498	7.634	ng	95
80) Benzo(a)anthracene	14.06	228	227211	7.727	ng	99
81) 3,3'-Dichlorobenzidine	14.02	252	78501	7.481	ng	98
82) Chrysene	14.10	228	225718	7.906	ng	97
83) Bis(2-ethylhexyl)phthalate	14.03	149	202771	8.049	ng #	99
84) Di-n-octyl phthalate	14.69	149	317108	7.791	ng	98
85) Indeno(1,2,3-cd)pyrene	17.12	276	194508	8.606	ng	97
87) Benzo(b)fluoranthene	15.16	252	207705	7.291	ng	98
88) Benzo(k)fluoranthene	15.19	252	218407	8.142	ng	99
89) Benzo(a)pyrene	15.54	252	195945	7.702	ng #	97
90) Dibenzo(a,h)anthracene	17.13	278	157016	7.987	ng	99
91) Benzo(g,h,i)perylene	17.58	276	161284	8.395	ng	98

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

