

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012618\
 Data File : BM013874.D
 Acq On : 27 Jan 2018 09:08
 Operator : SJ/JU
 Sample : J1222-13DL2 30X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6A95DL2

Manual Integrations
 APPROVED

Sohil
 1/29/2018 7:21:29 PM

Quant Time: Jan 27 22:48:28 2018
 Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM011018.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jan 27 04:13:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	136661	20.00	ng/ul	0.00
18) Naphthalene-d8	10.36	136	502798	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.23	164	271640	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.99	188	595063	20.00	ng/ul	0.00
75) Chrysene-d12	21.19	240	641420	20.00	ng/ul	0.00
83) Perylene-d12	23.38	264	664409	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.14	96	127	0.04	ng/uL	0.00
5) Phenol-d5	6.80	99	3228	0.31	ng/ul	0.03
7) Bis-(2-Chloroethyl)ether-d	6.95	67	3611	0.57	ng/ul	0.01
9) 2-Chlorophenol-d4	7.13	132	5061	0.59	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	3409m	0.42	ng/ul	0.04
19) Nitrobenzene-d5	8.77	128	1739m	0.45	ng/ul	0.04
22) 2-Nitrophenol-d4	9.48	143	1598	0.39	ng/ul	0.02
26) 2,4-Dichlorophenol-d3	10.06	165	1901	0.24	ng/ul	0.06
29) 4-Chloroaniline-d4	10.48	131	430	0.06	ng/ul	-0.03
43) Dimethylphthalate-d6	13.67	166	17189	0.77	ng/ul	0.01
46) Acenaphthylene-d8	13.93	160	20994	0.74	ng/ul	0.00
51) 4-Nitrophenol-d4	14.47	143	198	0.05	ng/ul	0.00
57) Fluorene-d10	15.24	176	17077	0.87	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.40	200	1250	0.35	ng/ul	0.02
70) Anthracene-d10	17.09	188	21157	0.73	ng/ul	0.00
76) Pyrene-d10	19.39	212	29356	0.98	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.24	264	26529	0.87	ng/ul	0.00

Target Compounds

					Qvalue	
49) Acenaphthene	14.30	153	68238	3.772	ng/ul	94
53) Dibenzofuran	14.65	168	56167	2.064	ng/ul	99
55) 2,3,4,6-Tetrachlorophenol	14.89	232	28202	4.946	ng/ul#	93
58) Fluorene	15.29	166	48547	2.178	ng/ul	93
68) Pentachlorophenol	16.64	266	573954	142.769	ng/ul	97
69) Phenanthrene	17.03	178	283699	8.577	ng/ul	98
71) Anthracene	17.13	178	60750m	1.781	ng/ul	
74) Fluoranthene	19.06	202	391654	9.919	ng/ul#	94
77) Pyrene	19.42	202	290324	7.445	ng/ul#	95
80) Benzo(a)anthracene	21.17	228	57591	1.486	ng/ul	98
82) Chrysene	21.20	228	93877m	2.640	ng/ul	
85) Benzo(b)fluoranthene	22.73	252	167245	4.099	ng/ul#	94
88) Benzo(a)pyrene	23.29	252	56469	1.465	ng/ul#	91

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

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