

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
 Data File : BM014397.D
 Acq On : 28 Feb 2018 11:00
 Operator : SJ/JU
 Sample : J1646-10DL 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE5X4DL

Manual Integrations
 APPROVED

Sohil
 3/1/2018 5:13:42 PM

Quant Time: Mar 01 10:10:34 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 28 12:00:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	101374	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	426808	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.17	164	238290	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.94	188	500621	20.00	ng/ul	0.00
75) Chrysene-d12	21.15	240	512014	20.00	ng/ul	0.00
83) Perylene-d12	23.33	264	488077	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.10	96	5738	2.42	ng/uL	0.00
5) Phenol-d5	6.72	99	124479	14.53	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	71841	13.85	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	97509	14.84	ng/ul	0.00
13) 4-Methylphenol-d8	8.25	113	100584	13.42	ng/ul	0.00
19) Nitrobenzene-d5	8.68	128	51556	16.35	ng/ul	0.00
22) 2-Nitrophenol-d4	9.40	143	56728	17.38	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.94	165	99071	14.65	ng/ul	0.01
29) 4-Chloroaniline-d4	10.46	131	108918m	16.35	ng/ul	0.00
43) Dimethylphthalate-d6	13.60	166	327556	16.65	ng/ul	0.00
46) Acenaphthylene-d8	13.87	160	411991	16.98	ng/ul	0.00
51) 4-Nitrophenol-d4	14.48	143	27434m	10.10	ng/ul	0.05
57) Fluorene-d10	15.18	176	276195	15.67	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.35	200	18925	6.30	ng/ul	0.01
70) Anthracene-d10	17.04	188	411585	17.04	ng/ul	0.00
76) Pyrene-d10	19.34	212	464191	17.58	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.19	264	398530	16.78	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.75	94	12934	1.439 ng/ul#	80
28) Naphthalene	10.34	128	24981	1.140 ng/ul#	95
44) Dimethylphthalate	13.65	163	79622	4.103 ng/ul	98
47) Acenaphthylene	13.90	152	70662	3.081 ng/ul#	94
49) Acenaphthene	14.24	153	39105	2.421 ng/ul	96
53) Dibenzofuran	14.59	168	68292m	2.820 ng/ul	
58) Fluorene	15.24	166	55892m	2.674 ng/ul	
69) Phenanthrene	16.98	178	1132567	39.266 ng/ul	99
71) Anthracene	17.07	178	187421	6.461 ng/ul	98
72) Carbazole	17.36	167	49862	2.044 ng/ul#	94
74) Fluoranthene	19.01	202	1224062	35.521 ng/ul#	92
77) Pyrene	19.37	202	1067890	31.384 ng/ul#	93
80) Benzo(a)anthracene	21.13	228	418088	13.084 ng/ul	99
81) Bis(2-ethylhexyl)phthalate	21.07	149	28537	1.416 ng/ul	99
82) Chrysene	21.19	228	373294	12.523 ng/ul	98
85) Benzo(b)fluoranthene	22.69	252	420707	12.873 ng/ul#	96
86) Benzo(k)fluoranthene	22.72	252	135327m	4.355 ng/ul	
88) Benzo(a)pyrene	23.23	252	300503	10.290 ng/ul#	94
89) Indeno(1,2,3-cd)pyrene	25.49	276	183151	6.463 ng/ul#	96
90) Dibenzo(a,h)anthracene	25.49	278	49355	2.096 ng/ul#	93
91) Benzo(a,h,i)perylene	26.14	276	130712	5.694 ng/ul	97

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

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