

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM121117\
 Data File : BM013321.D
 Acq On : 12 Dec 2017 01:21
 Operator : SJ/JU
 Sample : I6758-08 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9B09

Manual Integrations
 APPROVED

Sohil
 12/12/2017 3:57:12 PM

Quant Time: Dec 12 04:02:09 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 12 02:07:09 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	167611	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	756008	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.33	164	463884	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.08	188	1117270	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	1393899	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	1378706	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	2059	0.61	ng/uL	0.00
5) Phenol-d5	6.87	99	44797	3.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	26684	3.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	34072	3.01	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	38250	3.09	ng/ul	0.00
19) Nitrobenzene-d5	8.84	128	19022	3.17	ng/ul	0.00
22) 2-Nitrophenol-d4	9.56	143	20837	3.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	38683	2.93	ng/ul	0.00
29) 4-Chloroaniline-d4	10.61	131	35795	2.95	ng/ul	0.00
43) Dimethylphthalate-d6	13.75	166	140214	3.40	ng/ul	0.00
46) Acenaphthylene-d8	14.02	160	167962	3.30	ng/ul	0.00
51) 4-Nitrophenol-d4	14.55	143	19085m	2.66	ng/ul	0.01
57) Fluorene-d10	15.32	176	126348	3.56	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.45	200	14456	2.06	ng/ul	0.00
70) Anthracene-d10	17.17	188	202300	3.62	ng/ul	0.00
76) Pyrene-d10	19.47	212	227826	3.28	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.36	264	223028	3.18	ng/ul	0.00

Target Compounds

					Ovalue	
47) Acenaphthylene	14.05	152	458203	9.720	ng/ul	98
69) Phenanthrene	17.12	178	140202	2.209	ng/ul	97
71) Anthracene	17.21	178	436564	6.729	ng/ul	99
72) Carbazole	17.48	167	118383	2.121	ng/ul	97
74) Fluoranthene	19.14	202	1909358	24.546	ng/ul#	92
77) Pyrene	19.50	202	2195900	25.931	ng/ul#	92
80) Benzo(a)anthracene	21.26	228	2064843	23.348	ng/ul	94
82) Chrysene	21.31	228	3097763	37.755	ng/ul	98
85) Benzo(b)fluoranthene	22.84	252	5839314	62.828	ng/ul#	98
86) Benzo(k)fluoranthene	22.87	252	1479102m	16.911	ng/ul	
88) Benzo(a)pyrene	23.41	252	2612387	31.290	ng/ul#	98
89) Indeno(1,2,3-cd)pyrene	25.74	276	2130593	25.720	ng/ul#	97
90) Dibenzo(a,h)anthracene	25.74	278	592598	8.401	ng/ul#	88
91) Benzo(g,h,i)perylene	26.43	276	1768823	27.070	ng/ul#	95

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

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM121117\
Data File : BM013321.D
Acq On : 12 Dec 2017 01:21
Operator : SJ/JU
Sample : I6758-08 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F9B09

Manual Integrations
APPROVED

Sohil
12/12/2017 3:57:12 PM

Quant Time: Dec 12 04:02:09 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 12 02:07:09 2017
Response via : Initial Calibration

