

Data Path : Z:\HPCHEM1\BNA M\Data\BM011416\
 Data File : BM004008.D
 Acq On : 14 Jan 2016 05:50
 Operator : SJ/UM
 Sample : H1098-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BC963

Quant Time: Jan 14 12:51:17 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 13 06:45:02 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	43523	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	211032	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	123900	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	272937	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	255493	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	211666	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	2789	2.87	ng/uL	0.00
5) Phenol-d5	6.84	99	67781	18.76	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	41699	20.41	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	58471	19.69	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	52444	17.91	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	30053	18.91	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	32656	18.74	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	56935	18.39	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	68280	16.71	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	193255	19.87	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	240918	20.33	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	20898	11.96	ng/ul	0.00
58) Fluorene-d10	15.32	176	173832	21.07	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	18515	12.03	ng/ul	0.00
71) Anthracene-d10	17.17	188	270698	21.44	ng/ul	0.00
79) Pyrene-d10	19.47	212	291723	22.05	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	230832	21.84	ng/ul	0.00

Target Compounds

						Ovalue
45) Dimethylphthalate	13.77	163	110166	11.11	ng/ul	99
70) Phenanthrene	17.11	178	108044	7.04	ng/ul	100
72) Anthracene	17.20	178	20383	1.29	ng/ul	99
77) Fluoranthene	19.14	202	212084	13.30	ng/ul	98
80) Pyrene	19.50	202	303757	17.51	ng/ul	98
83) Benzo(a)anthracene	21.26	228	116097	7.63	ng/ul	97
85) Chrysene	21.32	228	140835	9.74	ng/ul	97
88) Benzo(b)fluoranthene	22.84	252	113269	8.80	ng/ul	98
89) Benzo(k)fluoranthene	22.84	252	28681	2.34	ng/ul#	97
91) Benzo(a)pyrene	23.41	252	96444	7.90	ng/ul	99
92) Indeno(1,2,3-cd)pyrene	25.76	276	52472	3.89	ng/ul	95
93) Dibenzo(a,h)anthracene	25.76	278	17829	1.57	ng/ul#	80
94) Benzo(g,h,i)perylene	26.45	276	54938	4.85	ng/ul	98

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

Data Path : Z:\HPCHEM1\BNA M\Data\BM011416\
Data File : BM004008.D
Acq On : 14 Jan 2016 05:50
Operator : SJ/UM
Sample : H1098-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BC963

Quant Time: Jan 14 12:51:17 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 13 06:45:02 2016
Response via : Initial Calibration

