

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
 Data File : BM002183.D
 Acq On : 02 Aug 2015 23:42
 Operator : TP
 Sample : G3095-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C01R4

Manual Integrations
 APPROVED

MMDadoda
 8/3/2015 4:44:56 PM

Quant Time: Aug 03 03:29:21 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 03 02:05:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	91415	20.00	ng/ul	0.00
18) Naphthalene-d8	10.35	136	410088	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.23	164	251942	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.98	188	575768	20.00	ng/ul	0.00
78) Chrysene-d12	21.19	240	580756	20.00	ng/ul	0.00
86) Perylene-d12	23.37	264	481440	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.01	96	2672	1.51	ng/uL	0.00
5) Phenol-d5	6.75	99	74872	11.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.89	67	41186	11.40	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	63394	11.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.28	113	59159	11.39	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	31920	10.95	ng/ul	0.00
22) 2-Nitrophenol-d4	9.44	143	37593	11.61	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	69524	11.15	ng/ul	0.00
29) 4-Chloroaniline-d4	10.49	131	59370	8.66	ng/ul	0.00
44) Dimethylphthalate-d6	13.64	166	208803	11.31	ng/ul	0.00
47) Acenaphthylene-d8	13.92	160	231461	10.03	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	21009	6.82	ng/ul	0.01
58) Fluorene-d10	15.23	176	160993	9.91	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.37	200	5446	1.55	ng/ul	0.00
71) Anthracene-d10	17.08	188	247138	9.62	ng/ul	0.00
79) Pyrene-d10	19.39	212	262136	10.72	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.23	264	215220	9.22	ng/ul	0.00

Target Compounds

						Qvalue
45) Dimethylphthalate	13.69	163	74309	4.03	ng/ul	99
70) Phenanthrene	17.02	178	348114	11.67	ng/ul	100
72) Anthracene	17.11	178	71934	2.36	ng/ul	100
77) Fluoranthene	19.05	202	542835	15.76	ng/ul	99
80) Pyrene	19.42	202	586500	18.56	ng/ul	99
83) Benzo(a)anthracene	21.17	228	298312	9.27	ng/ul	97
85) Chrysene	21.22	228	298701	9.92	ng/ul	98
88) Benzo(b)fluoranthene	22.72	252	335270	12.39	ng/ul	96
89) Benzo(k)fluoranthene	22.76	252	84351m	3.23	ng/ul	
91) Benzo(a)pyrene	23.28	252	230381	8.82	ng/ul	96
92) Indeno(1,2,3-cd)pyrene	25.57	276	145546	4.84	ng/ul#	95
93) Dibenzo(a,h)anthracene	25.57	278	39881	1.55	ng/ul#	85
94) Benzo(g,h,i)perylene	26.25	276	140330	5.56	ng/ul	96

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

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002183.D
Acq On : 02 Aug 2015 23:42
Operator : TP
Sample : G3095-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C01R4

Manual Integrations
APPROVED

MMDadoda
8/3/2015 4:44:56 PM

Quant Time: Aug 03 03:29:21 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 03 02:05:53 2015
Response via : Initial Calibration

