

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080415\
 Data File : BM002221.D
 Acq On : 04 Aug 2015 22:05
 Operator : TP/UM
 Sample : G3073-02RE
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 C01L3RE

Manual Integrations
 APPROVED

MMDadoda
 8/7/2015 4:42:43 PM

Quant Time: Aug 05 04:26:31 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 05 02:56:04 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	67049	20.00	ng/ul	0.00
18) Naphthalene-d8	10.31	136	275266	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.20	164	162818	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.95	188	363504	20.00	ng/ul	0.00
78) Chrysene-d12	21.17	240	403598	20.00	ng/ul	0.00
86) Perylene-d12	23.35	264	379833	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.97	96	3498	2.96	ng/uL	0.00
5) Phenol-d5	6.72	99	77703	15.60	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.86	67	41311	15.87	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	67332	15.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.25	113	63194	15.28	ng/ul	0.00
19) Nitrobenzene-d5	8.69	128	31495	15.78	ng/ul	0.00
22) 2-Nitrophenol-d4	9.40	143	36174	15.37	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.95	165	67067	14.89	ng/ul	0.00
29) 4-Chloroaniline-d4	10.46	131	63013	12.64	ng/ul	0.00
44) Dimethylphthalate-d6	13.61	166	187046	15.79	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	228196	15.25	ng/ul	0.00
52) 4-Nitrophenol-d4	14.45	143	18943	12.35	ng/ul	0.00
58) Fluorene-d10	15.20	176	156711	14.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.35	200	13291	6.35	ng/ul	0.00
71) Anthracene-d10	17.05	188	234208	14.48	ng/ul	0.00
79) Pyrene-d10	19.36	212	247493	12.81	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.21	264	272112	14.75	ng/ul	0.00

Target Compounds

					Qvalue
45) Dimethylphthalate	13.66	163	53493	4.53 ng/ul	99
48) Acenaphthylene	13.92	152	119624	7.84 ng/ul	100
59) Fluorene	15.25	166	20794	1.75 ng/ul#	96
70) Phenanthrene	17.00	178	323963	17.36 ng/ul	100
72) Anthracene	17.09	178	127173	6.67 ng/ul	98
77) Fluoranthene	19.03	202	1354040	66.29 ng/ul	99
80) Pyrene	19.39	202	1081638	43.60 ng/ul	99
83) Benzo(a)anthracene	21.15	228	755728	33.42 ng/ul	94
85) Chrysene	21.20	228	701787	33.71 ng/ul	98
88) Benzo(b)fluoranthene	22.70	252	1064587	49.74 ng/ul	98
89) Benzo(k)fluoranthene	22.74	252	331614m	16.00 ng/ul	
91) Benzo(a)pyrene	23.26	252	693709	33.94 ng/ul	98
92) Indeno(1,2,3-cd)pyrene	25.54	276	492053	21.36 ng/ul	97
93) Dibenzo(a,h)anthracene	25.54	278	140156	7.14 ng/ul#	82
94) Benzo(g,h,i)perylene	26.22	276	414752	21.62 ng/ul	100

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

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080415\
Data File : BM002221.D
Acq On : 04 Aug 2015 22:05
Operator : TP/UM
Sample : G3073-02RE
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C01L3RE

Quant Time: Aug 05 04:26:31 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080415.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 05 02:56:04 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

MMDadoda
8/7/2015 4:42:43 PM

