Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072216\
Data File : BM006635.D
Acq On : 22 Jul 2016 21:46
Operator : UM/SJ
Sample : H4077-15
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9YS9

Manual Integrations APPROVED
sohil
7/25/2016 6:39:08 PM

Quant Time: Jul 23 01:07:05 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 23 00:16:38 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	219846	20.00	ng/ul	0.00
7) Naphthalene-d8	10.40	136	971274	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.27	164	533819	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.03	188	1133336	20.00	ng/ul	0.01
29) Chrysene-d12	21.23	240	1119257	20.00	ng/ul	0.00
34) Perylene-d12	23.43	264	1202639	20.00	ng/ul	0.01
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.14	96	10995	2.02	ng/uL	0.00
3) Phenol-d5	6.80	99	424875	22.72	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.96	67	282801	24.41	ng/ul	0.00
5) 2-Chlorophenol-d4	7.16	132	344338	23.04	ng/ul	0.00
6) 4-Methylphenol-d8	8.33	113	301626	20.79	ng/ul	0.00
8) Nitrobenzene-d5	8.77	128	173291	23.57	ng/ul	0.00
9) 2-Nitrophenol-d4	9.49	143	177971	22.76	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.03	165	323624	21.90	ng/ul	0.01
12) 4-Chloroaniline-d4	10.55	131	357097	21.95	ng/ul	0.01
16) Dimethylphthalate-d6	13.69	166	944453	22.56	ng/ul	0.00
17) Acenaphthylene-d8	13.96	160	1230505	23.03	ng/ul	0.00
20) 4-Nitrophenol-d4	14.51	143	121811	16.76	ng/ul	0.02
21) Fluorene-d10	15.27	176	848095	22.66	ng/ul	0.01
24) 4,6-Dinitro-2-methylphenol	15.42	200	22376	3.80	ng/ul	0.01
26) Anthracene-d10	17.13	188	1250985	23.25	ng/ul	0.01
30) Pyrene-d10	19.43	212	1328855	26.06	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.29	264	1307016	23.63	ng/ul	0.01
Target Compounds						Qvalue
18) Acenaphthylene	13.99	152	97954	1.71	ng/ul	97
25) Phenanthrene	17.07	178	638770	10.02	ng/ul	98
27) Anthracene	17.16	178	126816	1.93	ng/ul	98
28) Fluoranthene	19.09	202	1713470	22.88	ng/ul	99
31) Pyrene	19.46	202	1401656	20.68	ng/ul	97
32) Benzo(a)anthracene	21.21	228	764151	11.27	ng/ul	96
33) Chrysene	21.26	228	861318	13.77	ng/ul	97
35) Benzo(b)fluoranthene	22.77	252	1296635	17.99	ng/ul	99
36) Benzo(k)fluoranthene	22.82	252	410521m	5.89	ng/ul	
38) Benzo(a)pyrene	23.34	252	806774	11.50	ng/ul	97
39) Indeno(1,2,3-cd)pyrene	25.64	276	637641	7.85	ng/ul	98
40) Dibenzo(a,h)anthracene	25.64	278	173233m	2.57	ng/ul	
41) Benzo(g,h,i)perylene	26.33	276	581658	8.35	ng/ul	98

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

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072216\
Data File : BM006635.D
Acq On : 22 Jul 2016 21:46
Operator : UM/SJ
Sample : H4077-15
Misc :
ALS Vial : 18 Sample Multiplier: 1

Quant Time: Jul 23 01:07:05 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 23 00:16:38 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
D9YS9

Manual Integrations
APPROVED
sohil
7/25/2016 6:39:08 PM