Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062716\
Data File : BM006070.D
Acq On : 27 Jun 2016 14:45
Operator : UM/SJ
Sample : H3669-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9Y61

Manual Integrations APPROVED
umangi
6/28/2016 2:34:20 PM

Quant Time: Jun 28 03:07:04 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 28 02:11:41 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	173677	20.00	ng/ul	0.00
7) Naphthalene-d8	10.44	136	911141	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.31	164	625744	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.06	188	1624562	20.00	ng/ul	0.00
29) Chrysene-d12	21.25	240	1905375	20.00	ng/ul	0.00
34) Perylene-d12	23.47	264	1727285	20.00	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	5359	1.38	ng/uL	0.00
3) Phenol-d5	6.83	99	278475	15.24	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.00	67	166886	16.42	ng/ul	0.00
5) 2-Chlorophenol-d4	7.19	132	210552	16.55	ng/ul	0.00
6) 4-Methylphenol-d8	8.36	113	177647	11.52	ng/ul	0.00
8) Nitrobenzene-d5	8.81	128	111595	16.09	ng/ul	0.00
9) 2-Nitrophenol-d4	9.53	143	116260	15.06	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.06	165	254986	17.22	ng/ul	0.00
12) 4-Chloroaniline-d4	10.58	131	243265	15.14	ng/ul	0.00
16) Dimethylphthalate-d6	13.72	166	919020	17.16	ng/ul	0.00
17) Acenaphthylene-d8	14.00	160	1162268	18.71	ng/ul	0.00
20) 4-Nitrophenol-d4	14.52	143	130012	11.38	ng/ul	-0.01
21) Fluorene-d10	15.30	176	861889	18.84	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.43	200	71803	7.96	ng/ul	-0.01
26) Anthracene-d10	17.16	188	1430731	19.17	ng/ul	0.00
30) Pyrene-d10	19.46	212	1756083	21.04	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.33	264	1514436	19.42	ng/ul	0.00
Target Compounds						Qvalue
25) Phenanthrene	17.10	178	189524	2.17	ng/ul	99
28) Fluoranthene	19.12	202	652439	5.83	ng/ul	99
31) Pyrene	19.49	202	557405	5.11	ng/ul	97
32) Benzo(a)anthracene	21.24	228	362725	3.20	ng/ul	97
33) Chrysene	21.29	228	427053	4.13	ng/ul	98
35) Benzo(b)fluoranthene	22.81	252	598232	5.96	ng/ul	97
36) Benzo(k)fluoranthene	22.84	252	170060m	1.82	ng/ul	
38) Benzo(a)pyrene	23.37	252	350406	3.64	ng/ul	98
39) Indeno(1,2,3-cd)pyrene	25.69	276	246700	2.15	ng/ul#	94
41) Benzo(g,h,i)perylene	26.37	276	220839	2.23	ng/ul	98

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

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062716\
Data File : BM006070.D
Acq On : 27 Jun 2016 14:45
Operator : UM/SJ
Sample : H3669-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Jun 28 03:07:04 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 28 02:11:41 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
D9Y61

Manual Integrations APPROVED
umangi
6/28/2016 2:34:20 PM