Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061816\
Data File : BM005870.D
Acq On : 17 Jun 2016 20:35
Operator : UM/SJ
Sample : H3537-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9Y43

Manual Integrations APPROVED
umangi
6/19/2016 12:14:53 AM

Quant Time: Jun 17 23:27:40 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061716.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 10:48:37 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	17722	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	79945	20.00	ng/ul	0.01
15) Acenaphthene-d10	14.39	164	50626	20.00	ng/ul	0.01
23) Phenanthrene-d10	17.13	188	118810	20.00	ng/ul	0.01
29) Chrysene-d12	21.32	240	120520	20.00	ng/ul	0.01
34) Perylene-d12	23.58	264	106126	20.00	ng/ul	0.02
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.18	96	449	3.28	ng/uL	0.00
3) Phenol-d5	6.93	99	29822	20.10	ng/ul	0.02
4) Bis-(2-Chloroethyl)ether-d	7.07	67	16333	21.24	ng/ul	0.00
5) 2-Chlorophenol-d4	7.27	132	28063	22.68	ng/ul	0.01
6) 4-Methylphenol-d8	8.46	113	23513	17.73	ng/ul	0.01
8) Nitrobenzene-d5	8.90	128	13944	24.71	ng/ul	0.00
9) 2-Nitrophenol-d4	9.62	143	16611	24.09	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	32506	23.87	ng/ul	0.01
12) 4-Chloroaniline-d4	10.67	131	20721m	22.48	ng/ul	0.00
16) Dimethylphthalate-d6	13.79	166	117043	24.44	ng/ul	0.00
17) Acenaphthylene-d8	14.07	160	140606	25.48	ng/ul	0.00
20) 4-Nitrophenol-d4	14.66	143	6416	11.12	ng/ul	0.04
21) Fluorene-d10	15.38	176	101662	24.91	ng/ul	0.01
24) 4,6-Dinitro-2-methylphenol	15.54	200	6926	8.64	ng/ul	0.02
26) Anthracene-d10	17.23	188	156407	25.21	ng/ul	0.01
30) Pyrene-d10	19.53	212	171801	30.07	ng/ul	0.01
37) Benzo(a)pyrene-d12	23.43	264	138556	26.11	ng/ul	0.02
Target Compounds						Qvalue
25) Phenanthrene	17.17	178	50999	7.21	ng/ul	99
27) Anthracene	17.26	178	8411m	1.16	ng/ul	
28) Fluoranthene	19.19	202	166311	18.71	ng/ul	99
31) Pyrene	19.56	202	239132	33.57	ng/ul	99
32) Benzo(a)anthracene	21.30	228	107516	14.28	ng/ul	97
33) Chrysene	21.36	228	135049	18.63	ng/ul	95
35) Benzo(b)fluoranthene	22.90	252	109726	16.61	ng/ul#	97
36) Benzo(k)fluoranthene	22.94	252	29548m	4.63	ng/ul	
38) Benzo(a)pyrene	23.48	252	70603	10.89	ng/ul	99
39) Indeno(1,2,3-cd)pyrene	25.87	276	39834	5.02	ng/ul	99
40) Dibenzo(a,h)anthracene	25.87	278	13803m	2.12	ng/ul	
41) Benzo(g,h,i)perylene	26.58	276	38266	5.67	ng/ul	97

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

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061816\
Data File : BM005870.D
Acq On : 17 Jun 2016 20:35
Operator : UM/SJ
Sample : H3537-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Quant Time: Jun 17 23:27:40 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061716.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 10:48:37 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
D9Y43

Manual Integrations APPROVED
umangi
6/19/2016 12:14:53 AM