Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM060716\
Data File : BM005700.D
Acq On : 07 Jun 2016 19:49
Operator : UM/SJ
Sample : H3411-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
D9XR7

Manual Integrations APPROVED
sohil
6/8/2016 7:16:50 PM

Quant Time: Jun 08 01:51:19 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM060216.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 08 01:24:32 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	117591	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	567952	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	322948	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	708161	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	634699	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	514623	20.00	ng/ul	-0.01
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.20	96	3772	1.35	ng/uL	0.00
3) Phenol-d5	6.93	99	146015	14.12	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.09	67	112504	17.88	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	134518	15.77	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	88081	10.77	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	76263	17.49	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	84883	18.19	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	133105	16.05	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	144184	14.63	ng/ul	0.00
16) Dimethylphthalate-d6	13.80	166	455535	18.35	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	590888	18.30	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	31199	8.78	ng/ul	0.00
21) Fluorene-d10	15.39	176	409387	19.20	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	35808	10.74	ng/ul	0.00
26) Anthracene-d10	17.24	188	626381	18.63	ng/ul	0.00
30) Pyrene-d10	19.54	212	661403	20.73	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	435399	18.38	ng/ul	0.00
Target Compounds						Qvalue
18) Acenaphthylene	14.12	152	35062	1.05	ng/ul	98
25) Phenanthrene	17.19	178	190731	4.88	ng/ul	98
27) Anthracene	17.27	178	41091	1.03	ng/ul	99
28) Fluoranthene	19.20	202	430437	10.77	ng/ul	99
31) Pyrene	19.57	202	385648	9.63	ng/ul	97
32) Benzo(a)anthracene	21.31	228	215730	5.80	ng/ul	97
33) Chrysene	21.36	228	224654	6.39	ng/ul	97
35) Benzo(b)fluoranthene	22.91	252	323177	10.41	ng/ul#	98
36) Benzo(k)fluoranthene	22.95	252	94583m	3.25	ng/ul	
38) Benzo(a)pyrene	23.49	252	183408	6.14	ng/ul#	95
39) Indeno(1,2,3-cd)pyrene	25.87	276	134940	3.82	ng/ul	95
40) Dibenzo(a,h)anthracene	25.87	278	37707	1.27	ng/ul#	79
41) Benzo(g,h,i)perylene	26.58	276	134329	4.53	ng/ul#	94

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

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM060716\
Data File : BM005700.D
Acq On : 07 Jun 2016 19:49
Operator : UM/SJ
Sample : H3411-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jun 08 01:51:19 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM060216.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 08 01:24:32 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
D9XR7

Manual Integrations APPROVED
sohil
6/8/2016 7:16:50 PM