Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011719\
Data File : BE098647.D
Acq On : 17 Jan 2019 9:53
Operator : SJ/JU
Sample : PB116433BS
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : PB116433BS

Quant Time: Jan 17 12:40:14 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE011519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 15 14:30:55 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	1107	0.40	ng		0.00
7) Naphthalene-d8	10.60	136	5209	0.40	ng		0.00
13) Acenaphthene-d10	14.47	164	3337	0.40	ng		0.00
19) Phenanthrene-d10	17.23	188	8116	0.40	ng		0.00
27) Chrysene-d12	21.41	240	9192	0.40	ng		0.00
34) Perylene-d12	23.92	264	9175	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.40	112	1161	0.39	ng		0.00
5) Phenol-d6	6.99	99	1710	0.41	ng		-0.01
8) Nitrobenzene-d5	8.96	82	1683	0.41	ng		0.00
11) 2-Methylnaphthalene-d10	12.21	152	3171	0.37	ng		0.00
14) 2,4,6-Tribromophenol	15.97	330	656	0.45	ng		0.00
15) 2-Fluorobiphenyl	13.09	172	5072	0.37	ng		0.00
25) Fluoranthene-d10	19.26	212	36609	0.36	ng		0.00
29) Terphenyl-d14	19.86	244	7728	0.38	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.31	88	732	0.352	ng		97
3) n-Nitrosodimethylamine	3.65	42	622	0.392	ng	#	98
6) bis(2-Chloroethyl)ether	7.23	93	1093	0.344	ng		95
9) Naphthalene	10.65	128	18839	0.363	ng		100
10) Hexachlorobutadiene	10.93	225	1430	0.385	ng		96
12) 2-Methylnaphthalene	12.28	142	3114	0.362	ng		95
16) Acenaphthylene	14.20	152	5489	0.380	ng		98
17) Acenaphthene	14.54	154	3474	0.383	ng		99
18) Fluorene	15.51	166	4475	0.379	ng		98
20) 4-Bromophenyl-phenylether	16.41	248	1593	0.371	ng		94
21) Hexachlorobenzene	16.53	284	1916	0.369	ng		94
22) Pentachlorophenol	16.89	266	693	0.526	ng		98
23) Phenanthrene	17.27	178	7114	0.370	ng		99
24) Anthracene	17.36	178	5828	0.354	ng		98
26) Fluoranthene	19.29	202	8983	0.363	ng		98
28) Pyrene	19.65	202	9263	0.366	ng		99
30) Benzo(a)anthracene	21.39	228	9631	0.390	ng		97
31) Chrysene	21.45	228	9380	0.366	ng		99
32) Bis(2-ethylhexyl)phthalate	21.32	149	18392	0.351	ng		100
33) Indeno(1,2,3-cd)pyrene	26.57	276	10293	0.390	ng		100
35) Benzo(b)fluoranthene	23.15	252	9130	0.365	ng		98
36) Benzo(k)fluoranthene	23.20	252	10158	0.371	ng		98
37) Benzo(a)pyrene	23.81	252	8962	0.365	ng		99
38) Dibenzo(a,h)anthracene	26.59	278	8518	0.391	ng	#	95
39) Benzo(g,h,i)perylene	27.37	276	8738	0.366	ng		98

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

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