Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE020419\
Data File : BE098713.D
Acq On : 4 Feb 2019 18:40
Operator : JU/SJ
Sample : K1308-01MSD
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SB-17(34.2-34.8)MSD

Manual Integrations APPROVED
apatel
2/5/2019 1:54:16 PM

Quant Time: Feb 04 23:58:57 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE012419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 24 14:44:44 2019
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	1108	0.40	ng	0.00
7) Naphthalene-d8	10.602	136	4105	0.40	ng	0.00
13) Acenaphthene-d10	14.471	164	2574	0.40	ng	0.00
19) Phenanthrene-d10	17.227	188	5926	0.40	ng	# 0.00
27) Chrysene-d12	21.415	240	7673	0.40	ng	# 0.00
34) Perylene-d12	23.920	264	7456	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.408	112	1028	0.31	ng	0.00
5) Phenol-d6	6.986	99	1567	0.33	ng	-0.01
8) Nitrobenzene-d5	8.978	82	1121	0.30	ng	0.00
11) 2-Methylnaphthalene-d10	12.210	152	2085m	0.27	ng	0.00
14) 2,4,6-Tribromophenol	15.969	330	388	0.34	ng	-0.01
15) 2-Fluorobiphenyl	13.088	172	2968	0.27	ng	-0.01
25) Fluoranthene-d10	19.259	212	25225	0.30	ng	0.00
29) Terphenyl-d14	19.856	244	4520	0.24	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.312	88	403	0.13	ng	# 44
3) n-Nitrosodimethylamine	3.646	42	916	0.36	ng	# 86
6) bis(2-Chloroethyl)ether	7.228	93	903	0.28	ng	96
9) Naphthalene	10.654	128	13115	0.28	ng	99
10) Hexachlorobutadiene	10.927	225	854	0.27	ng	96
12) 2-Methylnaphthalene	12.282	142	2078	0.27	ng	98
16) Acenaphthylene	14.198	152	3299	0.25	ng	99
17) Acenaphthene	14.529	154	2031	0.26	ng	99
18) Fluorene	15.514	166	2782	0.26	ng	99
20) 4-Bromophenyl-phenylether	16.410	248	977	0.28	ng	94
21) Hexachlorobenzene	16.531	284	1025	0.25	ng	# 91
22) Pentachlorophenol	16.892	266	577	0.55	ng	95
23) Phenanthrene	17.267	178	4699	0.29	ng	99
24) Anthracene	17.361	178	3701	0.29	ng	100
26) Fluoranthene	19.288	202	6055	0.29	ng	99
28) Pyrene	19.649	202	6151	0.25	ng	98
30) Benzo(a)anthracene	21.390	228	6546	0.28	ng	97
31) Chrysene	21.451	228	6401	0.26	ng	99
32) Bis(2-ethylhexyl)phtha...	21.318	149	16126	0.38	ng	99
33) Indeno(1,2,3-cd)pyrene	26.573	276	7171	0.27	ng	98
35) Benzo(b)fluoranthene	23.150	252	6580m	0.27	ng	
36) Benzo(k)fluoranthene	23.200	252	7446	0.29	ng	98
37) Benzo(a)pyrene	23.806	252	6475	0.28	ng	97
38) Dibenzo(a,h)anthracene	26.591	278	5902	0.28	ng	98
39) Benzo(g,h,i)perylene	27.375	276	6171	0.26	ng	99

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

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