Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE032219\
Data File : BE099080.D
Acq On : 22 Mar 2019 13:55
Operator : JU/SJ
Sample : SSTDICC3.2
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SSTDICC3.2

Manual Integrations APPROVED
Sohil
3/25/2019 12:11:45 PM

Quant Time: Mar 22 15:11:28 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 22 13:31:55 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	3128	0.40	ng	0.00
7) Naphthalene-d8	10.64	136	13005	0.40	ng	0.00
13) Acenaphthene-d10	14.49	164	9165	0.40	ng	0.00
19) Phenanthrene-d10	17.21	188	27086	0.40	ng	0.00
27) Chrysene-d12	21.39	240	33092	0.40	ng	0.00
34) Perylene-d12	23.90	264	31425	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.41	112	29664	3.17	ng	0.00
5) Phenol-d6	7.00	99	43415	3.29	ng	0.00
8) Nitrobenzene-d5	8.99	82	20226	1.48	ng	0.00
11) 2-Methylnaphthalene-d10	12.22	152	75942	3.16	ng	-0.01
14) 2,4,6-Tribromophenol	15.97	330	10980	2.37	ng	0.01
15) 2-Fluorobiphenyl	13.12	172	113827	3.00	ng	0.00
25) Fluoranthene-d10	19.24	212	1146871	2.65	ng	0.00
29) Terphenyl-d14	19.84	244	228870	2.85	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.35	88	15415	2.499	ng	93
3) n-Nitrosodimethylamine	3.66	42	9662m	1.228	ng	
6) bis(2-Chloroethyl)ether	7.27	93	37056	3.634	ng	98
9) Naphthalene	10.68	128	481683	3.234	ng	99
10) Hexachlorobutadiene	10.97	225	23311	2.366	ng	99
12) 2-Methylnaphthalene	12.31	142	84178	3.484	ng	99
16) Acenaphthylene	14.21	152	160604	3.407	ng	100
17) Acenaphthene	14.54	154	93719	3.352	ng	96
18) Fluorene	15.51	166	134551	3.337	ng	100
20) 4-Bromophenyl-phenylether	16.41	248	47558	3.240	ng	97
21) Hexachlorobenzene	16.52	284	44317	2.539	ng	98
22) Pentachlorophenol	16.87	266	15880	4.477	ng	98
23) Phenanthrene	17.25	178	252705	3.282	ng	99
24) Anthracene	17.35	178	249797	3.796	ng	100
26) Fluoranthene	19.28	202	353747	3.255	ng	99
28) Pyrene	19.63	202	361541	3.571	ng	100
30) Benzo(a)anthracene	21.37	228	368694	3.570	ng	99
31) Chrysene	21.43	228	365660	3.309	ng	100
32) Bis(2-ethylhexyl)phthalate	21.31	149	512023	2.330	ng	100
33) Indeno(1,2,3-cd)pyrene	26.56	276	396423	3.399	ng	99
35) Benzo(b)fluoranthene	23.13	252	346474	3.353	ng	98
36) Benzo(k)fluoranthene	23.18	252	350474	3.240	ng	96
37) Benzo(a)pyrene	23.79	252	327287	3.304	ng	96
38) Dibenzo(a,h)anthracene	26.59	278	325622	3.474	ng	98
39) Benzo(g,h,i)perylene	27.38	276	326143	3.327	ng	98

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

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