Quantitation Report (QT Reviewed)

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA E\DATA\BE090517\
Data File : BE093934.D
Acq On : 5 Sep 2017 13:12
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE090517

Manual Integrations APPROVED
Sohil
9/7/2017 6:30:54 PM

Quant Time: Sep 05 13:51:40 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 05 13:34:56 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1939	0.40	ng		0.00
7) Naphthalene-d8	10.71	136	10016	0.40	ng		0.00
13) Acenaphthene-d10	14.53	164	5748	0.40	ng		0.00
19) Phenanthrene-d10	17.27	188	9519	0.40	ng		0.00
27) Chrysene-d12	21.43	240	13689	0.40	ng		0.00
34) Perylene-d12	23.94	264	10892	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.53	112	2460	0.39	ng		0.00
5) Phenol-d6	7.11	99	3492	0.37	ng		0.00
8) Nitrobenzene-d5	9.08	82	3288	0.39	ng		0.00
11) 2-Methylnaphthalene-d10	12.30	152	6442	0.40	ng		0.00
14) 2,4,6-Tribromophenol	16.03	330	696	0.39	ng		0.00
15) 2-Fluorobiphenyl	13.16	172	9058	0.40	ng		0.00
25) Fluoranthene-d10	19.29	212	59597	0.42	ng		0.00
29) Terphenyl-d14	19.87	244	9872	0.39	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.43	88	1050	0.393	ng		99
3) n-Nitrosodimethylamine	3.75	42	1224	0.391	ng		97
6) bis(2-Chloroethyl)ether	7.34	93	2973	0.387	ng		86
9) Naphthalene	10.76	128	45344	0.395	ng		100
10) Hexachlorobutadiene	11.04	225	1599	0.383	ng		98
12) 2-Methylnaphthalene	12.37	142	7550	0.396	ng		97
16) Acenaphthylene	14.24	152	10868	0.379	ng		100
17) Acenaphthene	14.59	154	7605	0.392	ng		99
18) Fluorene	15.58	166	9661	0.386	ng		99
20) 4-Bromophenyl-phenylether	16.45	248	2511	0.470	ng		89
21) Hexachlorobenzene	16.58	284	2347	0.418	ng	#	100
22) Pentachlorophenol	16.94	266	608	0.403	ng		98
23) Phenanthrene	17.30	178	14249m	0.467	ng		
24) Anthracene	17.40	178	12702m	0.438	ng		
26) Fluoranthene	19.31	202	17699	0.417	ng		99
28) Pyrene	19.67	202	18327	0.389	ng		99
30) Benzo(a)anthracene	21.40	228	16499	0.392	ng		98
31) Chrysene	21.46	228	17571	0.395	ng		98
32) Bis(2-ethylhexyl)phthalate	21.30	149	39104	0.383	ng		99
33) Indeno(1,2,3-cd)pyrene	26.61	276	10977	0.370	ng		96
35) Benzo(b)fluoranthene	23.17	252	13107	0.378	ng		97
36) Benzo(k)fluoranthene	23.22	252	17568	0.413	ng		96
37) Benzo(a)pyrene	23.83	252	13988	0.396	ng		96
38) Dibenzo(a,h)anthracene	26.63	278	9221	0.394	ng	#	91
39) Benzo(g,h,i)perylene	27.42	276	9895	0.401	ng		98

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

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