Quantitation Report (QT/LSC Reviewed)

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA E\DATA\BE120517\
Data File : BE094943.D
Acq On : 6 Dec 2017 8:06
Operator : SJ/JU
Sample : I6707-03MSD
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : FT-MW01S-20171130MSD

Quant Time: Dec 06 10:01:38 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 01 13:33:28 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	7953	0.40	ng		-0.01
7) Naphthalene-d8	11.33	136	18113	0.40	ng		0.00
13) Acenaphthene-d10	15.06	164	10637	0.40	ng		-0.02
19) Phenanthrene-d10	17.77	188	25577	0.40	ng		0.00
27) Chrysene-d12	21.87	240	28423	0.40	ng		-0.02
34) Perylene-d12	24.55	264	22786	0.40	ng		-0.02
System Monitoring Compounds							
4) 2-Fluorophenol	5.96	112	2308	0.18	ng		0.00
5) Phenol-d6	7.59	99	2140	0.11	ng		0.00
8) Nitrobenzene-d5	9.66	82	6244	0.58	ng		0.00
11) 2-Methylnaphthalene-d10	12.86	152	12056	0.42	ng		0.00
14) 2,4,6-Tribromophenol	16.53	330	1205	0.52	ng		0.00
15) 2-Fluorobiphenyl	13.71	172	21034	0.46	ng		0.00
25) Fluoranthene-d10	19.75	212	140751	0.40	ng		0.00
29) Terphenyl-d14	20.31	244	28711	0.55	ng		-0.02
Target Compounds							Qvalue
2) 1,4-Dioxane	3.66	88	1227	0.147	ng	#	59
3) n-Nitrosodimethylamine	4.02	42	1750	0.165	ng	#	79
6) bis(2-Chloroethyl)ether	7.87	93	7357	0.398	ng		99
9) Naphthalene	11.38	128	87850	0.414	ng		99
10) Hexachlorobutadiene	11.64	225	3681	0.411	ng		97
12) 2-Methylnaphthalene	12.93	142	21508	0.408	ng		96
16) Acenaphthylene	14.79	152	22564	0.375	ng		99
17) Acenaphthene	15.12	154	14459	0.372	ng		96
18) Fluorene	16.10	166	18938	0.383	ng	#	78
20) 4-Bromophenyl-phenylether	16.97	248	5876	0.412	ng	#	85
21) Hexachlorobenzene	17.09	284	5645	0.406	ng	#	82
22) Pentachlorophenol	17.41	266	2639	0.706	ng	#	73
23) Phenanthrene	17.82	178	31894	0.391	ng		98
24) Anthracene	17.91	178	34201	0.396	ng	#	93
26) Fluoranthene	19.78	202	41046	0.392	ng		99
28) Pyrene	20.13	202	42553	0.434	ng		94
30) Benzo(a)anthracene	21.85	228	38097	0.437	ng	#	62
31) Chrysene	21.92	228	39362	0.409	ng	#	64
32) Bis(2-ethylhexyl)phthalate	21.75	149	90223	0.625	ng		100
33) Indeno(1,2,3-cd)pyrene	27.41	276	33152	0.421	ng		99
35) Benzo(b)fluoranthene	23.72	252	33111	0.398	ng	#	40
36) Benzo(k)fluoranthene	23.77	252	33549	0.396	ng	#	40
37) Benzo(a)pyrene	24.43	252	30480	0.418	ng	#	34
38) Dibenzo(a,h)anthracene	27.44	278	27508	0.408	ng	#	43
39) Benzo(g,h,i)perylene	28.30	276	28435	0.423	ng	#	53

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

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