Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062315\
Data File : BE089923.D
Acq On : 23 Jun 2015 20:48
Operator : TP/IZ
Sample : SSTDICC010
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC010

Quant Time: Jun 24 09:30:27 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 24 09:11:03 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	28430	5.00	ng		0.00
6) Naphthalene-d8	10.39	136	131408	5.00	ng		0.00
12) Acenaphthene-d10	14.25	164	80596	5.00	ng		0.00
18) Phenanthrene-d10	16.98	188	194034	5.00	ng		0.00
25) Chrysene-d12	21.14	240	243865	5.00	ng		0.00
32) Perylene-d12	23.47	264	213188	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.26	112	56249	8.77	ng		0.00
5) Phenol-d6	6.81	99	85659	10.18	ng		0.00
7) Nitrobenzene-d5	8.77	82	101200	10.03	ng		0.00
13) 2,4,6-Tribromophenol	15.74	330	14771	6.46	ng		0.00
14) 2-Fluorobiphenyl	12.87	172	222921	8.01	ng		0.00
27) Terphenyl-d14	19.60	244	398141	10.94	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.26	88	34108	10.57	ng		98
3) n-Nitrosodimethylamine	3.54	42	51957	10.50	ng	#	99
8) Nitrobenzene	8.82	77	108285	10.00	ng		97
9) Naphthalene	10.44	128	271768	8.62	ng		99
10) Hexachlorobutadiene	10.74	225	42193	7.98	ng		99
11) 2-Methylnaphthalene	12.05	142	191673	9.49	ng		99
15) Acenaphthylene	13.96	152	1392502	27.82	ng		99
16) Acenaphthene	14.31	154	199331	8.48	ng		97
17) Fluorene	15.29	166	261558	12.45	ng		98
19) 4-Bromophenyl-phenylether	16.20	248	76104	9.61	ng		95
20) Hexachlorobenzene	16.30	284	318735	34.60	ng		98
21) Pentachlorophenol	16.65	266	19874	8.82	ng		99
22) Phenanthrene	17.03	178	461367	7.20	ng		100
23) Anthracene	17.12	178	452638	15.25	ng		99
24) Fluoranthene	19.02	202	550446	12.79	ng		99
26) Pyrene	19.39	202	576424	7.01	ng		99
28) Benzo(a)anthracene	21.13	228	562924	7.90	ng		97
29) Chrysene	21.17	228	592969	7.92	ng		98
31) Indeno(1,2,3-cd)pyrene	25.90	276	553456	7.82	ng	#	100
33) Benzo(b)fluoranthene	22.76	252	501213	7.50	ng		99
34) Benzo(k)fluoranthene	22.81	252	533187	8.71	ng		99
35) Benzo(a)pyrene	23.37	252	462192	8.19	ng		100
36) Dibenzo(a,h)anthracene	25.92	278	449547	8.08	ng		100
37) Benzo(g,h,i)perylene	26.63	276	456770	7.79	ng		99

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

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