Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071615\
Data File : BE090140.D
Acq On : 17 Jul 2015 4:43
Operator : TP/IZ
Sample : G2948-09
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : YT-MW07

Quant Time: Jul 17 13:32:17 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 17 12:59:41 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	12157	5.00	ng		0.00
6) Naphthalene-d8	10.37	136	55346	5.00	ng		0.00
12) Acenaphthene-d10	14.22	164	31540	5.00	ng		0.00
18) Phenanthrene-d10	16.96	188	73929	5.00	ng		0.00
25) Chrysene-d12	21.12	240	76712	5.00	ng		0.00
32) Perylene-d12	23.44	264	63870	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.24	112	4941	1.90	ng		0.00
5) Phenol-d6	6.80	99	4610	1.25	ng		0.00
7) Nitrobenzene-d5	8.75	82	12658	3.09	ng		0.00
13) 2,4,6-Tribromophenol	15.73	330	4739	5.28	ng		0.00
14) 2-Fluorobiphenyl	12.84	172	29437	3.03	ng		0.00
27) Terphenyl-d14	19.58	244	55209	4.29	ng		0.00
Target Compounds							Qvalue
8) Nitrobenzene	8.79	77	317	0.07	ng		97
9) Naphthalene	10.41	128	1033	0.08	ng	#	83
10) Hexachlorobutadiene	10.71	225	112	0.05	ng		98
11) 2-Methylnaphthalene	12.03	142	756	0.09	ng	#	94
15) Acenaphthylene	13.93	152	5676	0.10	ng	#	99
16) Acenaphthene	14.29	154	817	0.10	ng		98
17) Fluorene	15.27	166	1265	0.12	ng	#	97
19) 4-Bromophenyl-phenylether	16.16	248	318	0.10	ng		83
20) Hexachlorobenzene	16.28	284	1372	0.10	ng	#	76
21) Pentachlorophenol	16.63	266	1576	2.04	ng		95
22) Phenanthrene	17.00	178	2011m	0.11	ng		
23) Anthracene	17.10	178	2036	0.12	ng		97
24) Fluoranthene	19.00	202	2212	0.12	ng		99
26) Pyrene	19.37	202	2368	0.11	ng		97
28) Benzo(a)anthracene	21.11	228	2075	0.16	ng	#	88
29) Chrysene	21.16	228	2237	0.04	ng	#	90
30) Bis(2-ethylhexyl)phthalate	21.05	149	3721	1.37	ng		98
31) Indeno(1,2,3-cd)pyrene	25.83	276	829	0.26	ng		98
33) Benzo(b)fluoranthene	22.73	252	885	0.19	ng	#	67
34) Benzo(k)fluoranthene	22.78	252	1532	0.25	ng	#	70
35) Benzo(a)pyrene	23.33	252	952	0.47	ng	#	59
36) Dibenzo(a,h)anthracene	25.85	278	466	0.17	ng	#	51
37) Benzo(g,h,i)perylene	26.57	276	943	0.44	ng	#	79

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

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