Quantitation Report (Not Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE063015\
Data File : BE090036.D
Acq On : 1 Jul 2015 7:46
Operator : TP/IZ
Sample : G2811-09
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : X1SB0510204-150625

Quant Time: Jul 01 17:22:44 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	16	5.00	ng		0.00
6) Naphthalene-d8	10.38	136	18	5.00	ng		0.00
12) Acenaphthene-d10	14.22	164	8	5.00	ng		-0.01
18) Phenanthrene-d10	16.96	188	55	5.00	ng		-0.02
25) Chrysene-d12	21.12	240	291	5.00	ng		-0.01
32) Perylene-d12	23.44	264	617	5.00	ng		-0.02
System Monitoring Compounds							
4) 2-Fluorophenol	5.26	112	12	3.26	ng		0.01
5) Phenol-d6	6.80	99	5	0.97	ng		0.00
7) Nitrobenzene-d5	8.76	82	1	0.70	ng		0.00
13) 2,4,6-Tribromophenol	15.74	330	1	4.27	ng		0.02
14) 2-Fluorobiphenyl	12.84	172	1	0.42	ng		-0.02
27) Terphenyl-d14	19.58	244	35	0.72	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.26	88	2	0.91	ng	#	1
3) n-Nitrosodimethylamine	3.53	42	10	3.36	ng	#	1
8) Nitrobenzene	8.80	77	2	1.32	ng	#	54
9) Naphthalene	10.41	128	3	0.74	ng	#	1
10) Hexachlorobutadiene	10.71	225	1	1.56	ng	#	1
11) 2-Methylnaphthalene	12.04	142	1	0.37	ng	#	50
15) Acenaphthylene	13.92	152	9	0.62	ng	#	1
16) Acenaphthene	14.32	154	2	0.96	ng	#	1
17) Fluorene	15.33	166	45	15.91	ng	#	46
19) 4-Bromophenyl-phenylether	16.13	248	6	2.57	ng	#	1
20) Hexachlorobenzene	16.28	284	17	1.75	ng	#	38
21) Pentachlorophenol	16.63	266	10	8.02	ng	#	58
22) Phenanthrene	17.01	178	45	3.19	ng	#	89
23) Anthracene	17.01	178	45	3.49	ng	#	88
24) Fluoranthene	19.01	202	112	7.12	ng	#	92
26) Pyrene	19.37	202	76	1.10	ng	#	62
28) Benzo(a)anthracene	21.11	228	407	5.60	ng	#	59
29) Chrysene	21.15	228	285	3.78	ng	#	59
30) Bis(2-ethylhexyl)phthalate	21.18	149	9	0.56	ng	#	1
31) Indeno(1,2,3-cd)pyrene	25.84	276	808	11.17	ng		95
33) Benzo(b)fluoranthene	22.73	252	1235	9.07	ng	#	47
34) Benzo(k)fluoranthene	22.73	252	1235	8.12	ng	#	47
35) Benzo(a)pyrene	23.33	252	748	5.95	ng	#	50
36) Dibenzo(a,h)anthracene	25.85	278	289	2.24	ng	#	18
37) Benzo(g,h,i)perylene	26.57	276	776	5.93	ng	#	60

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

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