

Data Path : S:\HPCHEM1\BNA_M\DATA\BM061115\
 Data File : BM001646.D
 Acq On : 11 Jun 2015 14:39
 Operator : TP/IZ
 Sample : G2414-08DL 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1T62DL

Quant Time: Jun 11 15:36:09 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 07:52:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1817	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	6523	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.34	164	2955	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.09	188	5674	0.40	ng/ul	0.00
18) Chrysene-d12	21.28	240	4254	0.40	ng/ul	0.01
22) Perylene-d12	23.52	264	3747	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	12.08	152	414	0.05	ng/ul	0.00
16) Fluoranthene-d10	19.12	212	766	0.07	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
5) Naphthalene	10.53	128	1748	0.10	ng/ul#	92
7) 2-Methylnaphthalene	12.15	142	1058	0.10	ng/ul	99
11) Fluorene	15.38	166	1513	0.13	ng/ul#	96
13) Pentachlorophenol	16.74	266	365	0.32	ng/ul	98
14) Phenanthrene	17.12	178	2469	0.15	ng/ul	96
17) Fluoranthene	19.14	202	2276	0.13	ng/ul	87
20) Benzo(a)anthracene	21.26	228	2267	0.17	ng/ul#	95
24) Benzo(k)fluoranthene	22.88	252	2509	0.17	ng/ul#	82
26) Indeno(1,2,3-cd)pyrene	25.77	276	2129	0.14	ng/ul	96
28) Benzo(g,h,i)perylene	26.45	276	2269	0.16	ng/ul#	89

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

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