

Data Path : S:\HPCHEM1\BNA_M\DATA\BM061115\
 Data File : BM001644.D
 Acq On : 11 Jun 2015 11:35
 Operator : TP/IZ
 Sample : G2414-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1T57

Quant Time: Jun 11 14:59:17 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	1645	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	6093	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.34	164	2785	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.09	188	5725	0.40	ng/ul	0.00
18) Chrysene-d12	21.27	240	3856	0.40	ng/ul	0.00
22) Perylene-d12	23.52	264	3037	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	2606	0.35	ng/ul	0.00
16) Fluoranthene-d10	19.12	212	4254	0.37	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
5) Naphthalene	10.53	128	5351	0.33	ng/ul	99
7) 2-Methylnaphthalene	12.15	142	5829	0.58	ng/ul	99
9) Acenaphthylene	14.05	152	3313	0.25	ng/ul	99
11) Fluorene	15.40	166	5738	0.53	ng/ul#	77
13) Pentachlorophenol	16.74	266	459	0.40	ng/ul	93
14) Phenanthrene	17.12	178	5451	0.33	ng/ul	99
15) Anthracene	17.22	178	8241	0.52	ng/ul	98
17) Fluoranthene	19.14	202	5265	0.30	ng/ul	96
19) Pyrene	19.52	202	6545	0.42	ng/ul	93
20) Benzo(a)anthracene	21.25	228	6200	0.50	ng/ul	96
21) Chrysene	21.31	228	6793	0.51	ng/ul	99
23) Benzo(b)fluoranthene	22.84	252	9033	0.77	ng/ul	96
24) Benzo(k)fluoranthene	22.88	252	4117	0.35	ng/ul	99
25) Benzo(a)pyrene	23.41	252	3709	0.34	ng/ul	98
27) Dibenzo(a,h)anthracene	25.79	278	6933	0.70	ng/ul	95
28) Benzo(g,h,i)perylene	26.45	276	3722	0.32	ng/ul	99

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

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