

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM071415\
 Data File : BM001915.D
 Acq On : 14 Jul 2015 14:15
 Operator : TP/UM
 Sample : G2928-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X1Y23

Quant Time: Jul 14 18:35:35 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 14 04:27:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1152	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	4570	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	2357	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	5133	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4428	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	3980	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	2140	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6142	0.44	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	5805	0.42	ng/ul	99
5) 2-Methylnaphthalene	12.12	142	4306	0.47	ng/ul	100
7) Acenaphthylene	14.03	152	7537	0.54	ng/ul#	90
8) Acenaphthene	14.37	153	5929	0.60	ng/ul	99
11) Pentachlorophenol	16.70	266	1560	1.15	ng/ul	98
13) Anthracene	17.18	178	9603	0.56	ng/ul	100
17) Pyrene	19.49	202	17112	0.85	ng/ul#	90
19) Chrysene	21.30	228	11900	0.69	ng/ul	91
21) Benzo(b)fluoranthene	22.81	252	13127	0.72	ng/ul	94
22) Benzo(k)fluoranthene	22.85	252	20116	1.33	ng/ul	94
23) Benzo(a)pyrene	23.38	252	10067	0.64	ng/ul	97
25) Dibenzo(a,h)anthracene	25.74	278	10485	0.78	ng/ul	98

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

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