

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM071415\
 Data File : BM001907.D
 Acq On : 13 Jul 2015 21:34
 Operator : TP/UM
 Sample : SSTD0.444
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.444

Quant Time: Jul 14 04:20:58 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 04:19:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1208	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	4646	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	2413	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	5307	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4712	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4298	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	2922	0.49	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	5821	0.49	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	5664	0.45	ng/ul	100
5) 2-Methylnaphthalene	12.12	142	3730	0.46	ng/ul	100
7) Acenaphthylene	14.01	152	5727	0.47	ng/ul	100
8) Acenaphthene	14.37	153	4034	0.44	ng/ul	100
9) Fluorene	15.37	166	4632	0.45	ng/ul	100
11) Pentachlorophenol	16.70	266	508	0.43	ng/ul	100
12) Phenanthrene	17.09	178	7447	0.45	ng/ul	100
13) Anthracene	17.18	178	7087	0.47	ng/ul	100
15) Fluoranthene	19.12	202	8230	0.46	ng/ul	100
17) Pyrene	19.48	202	8452	0.44	ng/ul	100
18) Benzo(a)anthracene	21.23	228	7309	0.46	ng/ul	100
19) Chrysene	21.28	228	7254	0.43	ng/ul	100
21) Benzo(b)fluoranthene	22.81	252	7648	0.44	ng/ul	100
22) Benzo(k)fluoranthene	22.85	252	6567	0.41	ng/ul	100
23) Benzo(a)pyrene	23.37	252	6695	0.44	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.71	276	7298	0.42	ng/ul	100
25) Dibenzo(a,h)anthracene	25.72	278	5852	0.41	ng/ul	100
26) Benzo(g,h,i)perylene	26.40	276	6254	0.41	ng/ul	100

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

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