

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061815\
 Data File : BM001727.D
 Acq On : 17 Jun 2015 20:39
 Operator : TP/IZ
 Sample : SSTD0.406
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.406

Quant Time: Jun 18 04:23:55 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 18 04:21:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	995	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	3582	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	1675	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	3542	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	2844	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	2494	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	1759	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	2987	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.50	128	3799	0.42	ng/ul	100
5) 2-Methylnaphthalene	12.12	142	2404	0.40	ng/ul	100
7) Acenaphthylene	14.03	152	3210	0.43	ng/ul	100
8) Acenaphthene	14.37	153	2463	0.42	ng/ul	100
9) Fluorene	15.37	166	2752	0.42	ng/ul	100
11) Pentachlorophenol	16.72	266	286	0.37	ng/ul	100
12) Phenanthrene	17.11	178	4197	0.41	ng/ul	100
13) Anthracene	17.18	178	3823	0.41	ng/ul	100
15) Fluoranthene	19.13	202	4484	0.40	ng/ul	100
17) Pyrene	19.49	202	4572	0.41	ng/ul	100
18) Benzo(a)anthracene	21.23	228	3715	0.42	ng/ul	100
19) Chrysene	21.28	228	3980	0.40	ng/ul	100
21) Benzo(b)fluoranthene	22.80	252	3741	0.41	ng/ul	100
22) Benzo(k)fluoranthene	22.85	252	3774	0.41	ng/ul	100
23) Benzo(a)pyrene	23.37	252	3436	0.40	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.70	276	3952	0.41	ng/ul	100
25) Dibenzo(a,h)anthracene	25.73	278	3202	0.42	ng/ul	100
26) Benzo(g,h,i)perylene	26.39	276	3501	0.41	ng/ul	100

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

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