

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020515\
 Data File : BM000425.D
 Acq On : 05 Feb 2015 11:42
 Operator : TP/IZ
 Sample : SSTD0.476
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.476

Quant Time: Feb 05 13:08:24 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM020515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 05 12:58:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2210	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7191	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4385	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	8988	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	7772	0.40	ng/ul	0.00
20) Perylene-d12	23.55	264	6808	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.11	152	3737	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	7544	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.56	128	7061	0.39	ng/ul	100
5) 2-Methylnaphthalene	12.18	142	4854	0.37	ng/ul	100
7) Acenaphthylene	14.10	152	6873	0.38	ng/ul	100
8) Acenaphthene	14.43	153	5231	0.39	ng/ul	100
9) Fluorene	15.43	166	6224	0.37	ng/ul	100
11) Pentachlorophenol	16.78	266	791	0.69	ng/ul	100
12) Phenanthrene	17.16	178	9868	0.38	ng/ul	100
13) Anthracene	17.26	178	8999	0.37	ng/ul	100
15) Fluoranthene	19.19	202	11096	0.37	ng/ul	100
17) Pyrene	19.55	202	11399	0.36	ng/ul	100
18) Benzo(a)anthracene	21.29	228	9237	0.39	ng/ul	100
19) Chrysene	21.34	228	10279	0.38	ng/ul	100
21) Benzo(b)fluoranthene	22.88	252	9127	0.36	ng/ul	100
22) Benzo(k)fluoranthene	22.93	252	9968	0.37	ng/ul	100
23) Benzo(a)pyrene	23.46	252	8747	0.37	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.82	276	9998	0.42	ng/ul	100
25) Dibenzo(a,h)anthracene	25.84	278	7939	0.42	ng/ul	100
26) Benzo(g,h,i)perylene	26.51	276	8736	0.42	ng/ul	100

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

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