

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021315\
 Data File : BM000520.D
 Acq On : 13 Feb 2015 15:24
 Operator : TP/IZ
 Sample : SSTD0.406
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.498

Quant Time: Feb 13 16:43:04 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 12 16:45:55 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	3415	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	11237	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	6420	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	13583	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	11871	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10226	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	5593	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	11330	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	10715	0.41	ng/ul	100
5) 2-Methylnaphthalene	12.16	142	7159	0.40	ng/ul	100
7) Acenaphthylene	14.08	152	9925	0.41	ng/ul	99
8) Acenaphthene	14.41	153	7790	0.42	ng/ul	98
9) Fluorene	15.40	166	9299	0.42	ng/ul	100
11) Pentachlorophenol	16.75	266	569	0.33	ng/ul	97
12) Phenanthrene	17.13	178	15068	0.41	ng/ul	98
13) Anthracene	17.23	178	13553	0.41	ng/ul	98
15) Fluoranthene	19.15	202	16975	0.41	ng/ul	92
17) Pyrene	19.51	202	17676	0.41	ng/ul	92
18) Benzo(a)anthracene	21.27	228	13411	0.40	ng/ul	98
19) Chrysene	21.32	228	15924	0.41	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	15251	0.44	ng/ul	96
22) Benzo(k)fluoranthene	22.89	252	13192	0.38	ng/ul	99
23) Benzo(a)pyrene	23.41	252	13096	0.41	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.75	276	13848	0.38	ng/ul	100
25) Dibenzo(a,h)anthracene	25.76	278	10751	0.38	ng/ul	97
26) Benzo(g,h,i)perylene	26.43	276	12379	0.39	ng/ul	99

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

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