

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021115\
 Data File : BM000498.D
 Acq On : 12 Feb 2015 04:04
 Operator : TP/IZ
 Sample : SSTD0.404
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.404

Quant Time: Feb 12 07:34:28 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 11 23:58:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	2971	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	10181	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6409	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14587	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	13370	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	11127	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.10	152	5334	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	12646	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.55	128	9694	0.39	ng/ul	99
5) 2-Methylnaphthalene	12.17	142	6713	0.39	ng/ul	99
7) Acenaphthylene	14.09	152	10418	0.41	ng/ul	99
8) Acenaphthene	14.41	153	7850	0.40	ng/ul	98
9) Fluorene	15.42	166	9563	0.40	ng/ul	99
11) Pentachlorophenol	16.76	266	696	0.52	ng/ul	98
12) Phenanthrene	17.15	178	15729	0.38	ng/ul	99
13) Anthracene	17.24	178	14457	0.39	ng/ul	99
15) Fluoranthene	19.16	202	18631	0.39	ng/ul	91
17) Pyrene	19.54	202	19179	0.39	ng/ul	100
18) Benzo(a)anthracene	21.28	228	15072	0.38	ng/ul	95
19) Chrysene	21.33	228	17444	0.39	ng/ul	98
21) Benzo(b)fluoranthene	22.86	252	15738	0.41	ng/ul	97
22) Benzo(k)fluoranthene	22.90	252	14834	0.36	ng/ul	98
23) Benzo(a)pyrene	23.43	252	14141	0.39	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.77	276	13831	0.36	ng/ul	98
25) Dibenzo(a,h)anthracene	25.79	278	10976	0.37	ng/ul	98
26) Benzo(g,h,i)perylene	26.46	276	11867	0.36	ng/ul	99

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

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