

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021315\
 Data File : BM000516.D
 Acq On : 13 Feb 2015 08:04
 Operator : TP/IZ
 Sample : SSTD0.406
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.406

Quant Time: Feb 13 12:03:15 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	3284	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	10968	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	6395	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	13707	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	11999	0.40	ng/ul	-0.01
20) Perylene-d12	23.51	264	10358	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	5471	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	11503	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	10430	0.41	ng/ul	99
5) 2-Methylnaphthalene	12.16	142	6963	0.40	ng/ul	100
7) Acenaphthylene	14.07	152	9825	0.41	ng/ul	99
8) Acenaphthene	14.40	153	7584	0.41	ng/ul#	79
9) Fluorene	15.40	166	9072	0.41	ng/ul	96
11) Pentachlorophenol	16.75	266	668	0.38	ng/ul	99
12) Phenanthrene	17.13	178	14597	0.39	ng/ul	99
13) Anthracene	17.23	178	13288	0.40	ng/ul	99
15) Fluoranthene	19.15	202	16834	0.41	ng/ul	91
17) Pyrene	19.51	202	17568	0.41	ng/ul	92
18) Benzo(a)anthracene	21.26	228	13668	0.40	ng/ul	92
19) Chrysene	21.31	228	15905	0.41	ng/ul	92
21) Benzo(b)fluoranthene	22.83	252	13829	0.40	ng/ul	98
22) Benzo(k)fluoranthene	22.88	252	14750	0.42	ng/ul	98
23) Benzo(a)pyrene	23.41	252	13274	0.41	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.74	276	13983	0.38	ng/ul	100
25) Dibenzo(a,h)anthracene	25.76	278	10888	0.38	ng/ul	98
26) Benzo(g,h,i)perylene	26.43	276	12330	0.38	ng/ul	99

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

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