

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120914\
 Data File : BE089203.D
 Acq On : 10 Dec 2014 1:27
 Operator : TP/JJ
 Sample : SSTD0.438
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.438

Quant Time: Dec 10 02:16:02 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM01.2-EPA-SIM-BE120514.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 10 01:02:19 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	2607	0.40	ng/ul	0.00
2) Naphthalene-d8	8.51	136	10259	0.40	ng/ul	0.00
6) Acenaphthene-d10	10.65	164	4942	0.40	ng/ul	0.00
10) Phenanthrene-d10	12.71	188	7054	0.40	ng/ul	0.00
16) Chrysene-d12	18.06	240	4725	0.40	ng/ul	0.00
20) Perylene-d12	21.11	264	4153	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	9.35	152	5036	0.42	ng/ul	0.00
14) Fluoranthene-d10	14.91	212	4917	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	8.53	128	10192	0.41	ng/ul	100
5) 2-Methylnaphthalene	9.38	142	6603	0.44	ng/ul	100
7) Acenaphthylene	10.48	152	36678	0.42	ng/ul	100
8) Acenaphthene	10.69	153	22092	0.42	ng/ul	93
9) Fluorene	11.34	166	6029	0.42	ng/ul	95
11) Pentachlorophenol	12.43	266	413	0.33	ng/ul	98
12) Phenanthrene	12.75	178	32061	0.42	ng/ul	98
13) Anthracene	12.83	178	31596	0.42	ng/ul	100
15) Fluoranthene	14.95	202	26893	0.39	ng/ul	99
17) Pyrene	15.40	202	27403	0.42	ng/ul	98
18) Benzo(a)anthracene	18.04	228	5189	0.43	ng/ul	99
19) Chrysene	18.12	228	5258	0.42	ng/ul	99
21) Benzo(b)fluoranthene	20.36	252	4449	0.43	ng/ul	99
22) Benzo(k)fluoranthene	20.42	252	5185	0.43	ng/ul	99
23) Benzo(a)pyrene	21.00	252	4693	0.43	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	23.21	276	14425	0.42	ng/ul	100
25) Dibenzo(a,h)anthracene	23.29	278	3829	0.43	ng/ul	99
26) Benzo(g,h,i)perylene	23.79	276	17362	0.41	ng/ul	100

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

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