

Data Path : Z:\HPCHEM1\BNA_E\Data\BE120814\
 Data File : BE089175.D
 Acq On : 8 Dec 2014 15:23
 Operator : TP/JJ
 Sample : SSTD0.436
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.436

Quant Time: Dec 09 00:42:30 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 : Tue Dec 09 00:37:23 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	2683	0.40	ng/ul	0.00
2) Naphthalene-d8	8.51	136	10562	0.40	ng/ul	0.00
6) Acenaphthene-d10	10.66	164	5279	0.40	ng/ul	0.00
10) Phenanthrene-d10	12.73	188	7755	0.40	ng/ul	0.00
16) Chrysene-d12	18.08	240	5328	0.40	ng/ul	0.00
20) Perylene-d12	21.13	264	4715	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	9.36	152	5170	0.42	ng/ul	0.00
14) Fluoranthene-d10	14.93	212	5665	0.41	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	8.54	128	10468	0.41	ng/ul	98
5) 2-Methylnaphthalene	9.39	142	6739	0.43	ng/ul	100
7) Acenaphthylene	10.48	152	38006	0.41	ng/ul	99
8) Acenaphthene	10.69	153	23226	0.41	ng/ul	89
9) Fluorene	11.35	166	6413	0.42	ng/ul	96
11) Pentachlorophenol	12.45	266	440	0.32	ng/ul	99
12) Phenanthrene	12.76	178	36003	0.43	ng/ul	99
13) Anthracene	12.85	178	34696	0.42	ng/ul	99
15) Fluoranthene	14.97	202	31048	0.41	ng/ul	98
17) Pyrene	15.42	202	31658	0.43	ng/ul	99
18) Benzo(a)anthracene	18.06	228	5991	0.44	ng/ul	99
19) Chrysene	18.14	228	5847	0.41	ng/ul	97
21) Benzo(b)fluoranthene	20.38	252	5312	0.46	ng/ul	99
22) Benzo(k)fluoranthene	20.43	252	5578	0.41	ng/ul	100
23) Benzo(a)pyrene	21.02	252	5351	0.44	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	23.22	276	18128	0.46	ng/ul	100
25) Dibenzo(a,h)anthracene	23.32	278	4530	0.45	ng/ul	99
26) Benzo(g,h,i)perylene	23.81	276	20382	0.42	ng/ul	100

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

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