

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120914\
 Data File : BE089197.D
 Acq On : 9 Dec 2014 21:31
 Operator : TP/JJ
 Sample : SSTD0.437
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.437

Quant Time: Dec 10 01:04:26 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 : Fri Dec 05 23:47:26 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	2519	0.40	ng/ul	-0.02
2) Naphthalene-d8	8.51	136	9966	0.40	ng/ul	-0.02
6) Acenaphthene-d10	10.65	164	4934	0.40	ng/ul	-0.03
10) Phenanthrene-d10	12.71	188	7211	0.40	ng/ul	-0.02
16) Chrysene-d12	18.06	240	4728	0.40	ng/ul	-0.03
20) Perylene-d12	21.11	264	3978	0.40	ng/ul	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	9.35	152	4902	0.42	ng/ul	-0.02
14) Fluoranthene-d10	14.91	212	5153	0.40	ng/ul	-0.03

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	8.54	128	9903	0.41	ng/ul	98
5) 2-Methylnaphthalene	9.39	142	6453	0.44	ng/ul	100
7) Acenaphthylene	10.47	152	35768	0.41	ng/ul	100
8) Acenaphthene	10.69	153	22148	0.42	ng/ul	97
9) Fluorene	11.34	166	6017	0.42	ng/ul	99
11) Pentachlorophenol	12.45	266	365	0.29	ng/ul	99
12) Phenanthrene	12.75	178	32955	0.42	ng/ul	99
13) Anthracene	12.83	178	31986	0.42	ng/ul	100
15) Fluoranthene	14.95	202	28173	0.40	ng/ul	99
17) Pyrene	15.40	202	28498	0.44	ng/ul	99
18) Benzo(a)anthracene	18.04	228	4977	0.41	ng/ul	99
19) Chrysene	18.12	228	5268	0.42	ng/ul	99
21) Benzo(b)fluoranthene	20.36	252	4145	0.42	ng/ul	99
22) Benzo(k)fluoranthene	20.42	252	4975	0.43	ng/ul	100
23) Benzo(a)pyrene	21.00	252	4312	0.42	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	23.20	276	13338	0.40	ng/ul	99
25) Dibenzo(a,h)anthracene	23.29	278	3578	0.42	ng/ul	99
26) Benzo(g,h,i)perylene	23.78	276	16637	0.41	ng/ul	100

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

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