

Data Path : Z:\HPCHEM1\BNA_E\Data\BE120814\
 Data File : BE089173.D
 Acq On : 8 Dec 2014 13:14
 Operator : TP/JJ
 Sample : SSTD0.435
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.435

Quant Time: Dec 09 00:39:23 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	2418	0.40	ng/ul	-0.01
2) Naphthalene-d8	8.51	136	9700	0.40	ng/ul	-0.02
6) Acenaphthene-d10	10.66	164	4835	0.40	ng/ul	-0.02
10) Phenanthrene-d10	12.73	188	7345	0.40	ng/ul	-0.01
16) Chrysene-d12	18.08	240	5257	0.40	ng/ul	0.00
20) Perylene-d12	21.14	264	4660	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	9.36	152	4772	0.42	ng/ul	-0.01
14) Fluoranthene-d10	14.93	212	5445	0.42	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	8.54	128	9766	0.42	ng/ul	99
5) 2-Methylnaphthalene	9.39	142	6216	0.43	ng/ul	100
7) Acenaphthylene	10.48	152	35339	0.42	ng/ul	99
8) Acenaphthene	10.70	153	21517	0.42	ng/ul	96
9) Fluorene	11.35	166	5965	0.42	ng/ul	98
11) Pentachlorophenol	12.45	266	431	0.33	ng/ul	97
12) Phenanthrene	12.76	178	32817	0.41	ng/ul	98
13) Anthracene	12.85	178	32610	0.42	ng/ul	100
15) Fluoranthene	14.97	202	29473	0.41	ng/ul	98
17) Pyrene	15.42	202	30498	0.42	ng/ul	98
18) Benzo(a)anthracene	18.06	228	5961	0.44	ng/ul	100
19) Chrysene	18.14	228	5733	0.41	ng/ul	99
21) Benzo(b)fluoranthene	20.38	252	5285	0.46	ng/ul	99
22) Benzo(k)fluoranthene	20.44	252	5453	0.40	ng/ul	99
23) Benzo(a)pyrene	21.02	252	5262	0.43	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	23.23	276	17927	0.46	ng/ul	99
25) Dibenzo(a,h)anthracene	23.32	278	4496	0.45	ng/ul	100
26) Benzo(g,h,i)perylene	23.82	276	19770	0.41	ng/ul	99

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

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