

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071415\
 Data File : BE090101.D
 Acq On : 14 Jul 2015 16:42
 Operator : TP/IZ
 Sample : SSTD0.804
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.804

Quant Time: Jul 15 00:46:09 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM02.2-EPA-SIM-BE071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 15 00:44:52 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	1488	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6978	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	4017	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	8780	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	9096	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	6305	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.96	152	9160	0.83	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	21528	0.89	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	15426	0.74	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	10410	0.75	ng/ul	100
7) Acenaphthylene	13.93	152	68125	1.90	ng/ul	99
8) Acenaphthene	14.28	153	47713	1.90	ng/ul	99
9) Fluorene	15.27	166	13950	0.74	ng/ul	98
11) Pentachlorophenol	16.61	266	636	0.31	ng/ul	94
12) Phenanthrene	16.99	178	88032	1.93	ng/ul	97
13) Anthracene	17.07	178	88251	2.01	ng/ul	96
15) Fluoranthene	19.00	202	98171	1.94	ng/ul	99
17) Pyrene	19.37	202	100353	1.72	ng/ul	100
18) Benzo(a)anthracene	21.11	228	20052	0.60	ng/ul	99
19) Chrysene	21.16	228	24789	0.72	ng/ul	99
21) Benzo(b)fluoranthene	22.73	252	15462	0.58	ng/ul	98
22) Benzo(k)fluoranthene	22.78	252	22292	0.91	ng/ul	98
23) Benzo(a)pyrene	23.33	252	16532	0.69	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.85	276	65753	1.78	ng/ul	99
25) Dibenzo(a,h)anthracene	25.86	278	12660	0.63	ng/ul	93
26) Benzo(g,h,i)perylene	26.58	276	61314	1.88	ng/ul	97

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

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