

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071415\
 Data File : BE090100.D
 Acq On : 14 Jul 2015 16:06
 Operator : TP/IZ
 Sample : SSTD0.403
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.403

Quant Time: Jul 15 00:45:56 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	1448	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6549	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	3777	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	8740	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	8642	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	6143	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.96	152	4256	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	10176	0.42	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	7202	0.37	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	4787	0.37	ng/ul	100
7) Acenaphthylene	13.93	152	31854	0.95	ng/ul	100
8) Acenaphthene	14.28	153	22099	0.94	ng/ul	100
9) Fluorene	15.27	166	6561	0.37	ng/ul	100
11) Pentachlorophenol	16.61	266	311	0.15	ng/ul	100
12) Phenanthrene	16.99	178	42236	0.93	ng/ul	100
13) Anthracene	17.08	178	40745	0.93	ng/ul	100
15) Fluoranthene	19.00	202	46534	0.92	ng/ul	100
17) Pyrene	19.37	202	47611	0.86	ng/ul	100
18) Benzo(a)anthracene	21.11	228	9072	0.29	ng/ul	100
19) Chrysene	21.16	228	11672	0.36	ng/ul	100
21) Benzo(b)fluoranthene	22.73	252	6758	0.26	ng/ul	100
22) Benzo(k)fluoranthene	22.78	252	10125	0.43	ng/ul	100
23) Benzo(a)pyrene	23.33	252	7581	0.32	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.85	276	30064	0.84	ng/ul	100
25) Dibenzo(a,h)anthracene	25.86	278	5610	0.29	ng/ul	100
26) Benzo(g,h,i)perylene	26.58	276	28298	0.89	ng/ul	100

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

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