

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071415\
 Data File : BE090102.D
 Acq On : 14 Jul 2015 17:18
 Operator : TP/IZ
 Sample : SSTD1.605
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.605

Quant Time: Jul 15 00:46:20 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	1463	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6795	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	3851	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	8504	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	9406	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	6623	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.96	152	17840	1.67	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	43059	1.84	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	29978	1.48	ng/ul	98
5) 2-Methylnaphthalene	12.03	142	20261	1.51	ng/ul	100
7) Acenaphthylene	13.93	152	133260	3.88	ng/ul	99
8) Acenaphthene	14.28	153	92082	3.82	ng/ul	98
9) Fluorene	15.27	166	26987	1.49	ng/ul	98
11) Pentachlorophenol	16.61	266	1412	0.71	ng/ul	95
12) Phenanthrene	16.99	178	169990	3.85	ng/ul	98
13) Anthracene	17.07	178	172773	4.06	ng/ul	96
15) Fluoranthene	19.00	202	196563	4.01	ng/ul	99
17) Pyrene	19.37	202	201279	3.35	ng/ul	100
18) Benzo(a)anthracene	21.11	228	42904	1.24	ng/ul	99
19) Chrysene	21.16	228	51164	1.43	ng/ul	99
21) Benzo(b)fluoranthene	22.73	252	34370	1.22	ng/ul	98
22) Benzo(k)fluoranthene	22.78	252	47844	1.87	ng/ul	97
23) Benzo(a)pyrene	23.33	252	36178	1.44	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.85	276	146582	3.78	ng/ul	99
25) Dibenzo(a,h)anthracene	25.86	278	28729	1.37	ng/ul	92
26) Benzo(g,h,i)perylene	26.58	276	134249	3.92	ng/ul	97

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

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