

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021115\
 Data File : BM000483.D
 Acq On : 11 Feb 2015 18:37
 Operator : TP/IZ
 Sample : SST00.403
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SST00.403

Quant Time: Feb 11 23:59:16 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 11 14:38:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	3093	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	10233	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6152	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13014	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	9563	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	7586	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.10	152	5223	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	10329	0.36	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.55	128	9683	0.39	ng/ul	99
5) 2-Methylnaphthalene	12.17	142	6668	0.38	ng/ul	99
7) Acenaphthylene	14.09	152	9515	0.39	ng/ul	99
8) Acenaphthene	14.41	153	7468	0.40	ng/ul	99
9) Fluorene	15.42	166	8884	0.39	ng/ul	100
11) Pentachlorophenol	16.76	266	314	0.26	ng/ul	93
12) Phenanthrene	17.15	178	13988	0.38	ng/ul	99
13) Anthracene	17.24	178	12655	0.38	ng/ul	99
15) Fluoranthene	19.16	202	15099	0.36	ng/ul	90
17) Pyrene	19.54	202	15155	0.43	ng/ul	100
18) Benzo(a)anthracene	21.28	228	10389	0.37	ng/ul	99
19) Chrysene	21.33	228	12834	0.40	ng/ul	99
21) Benzo(b)fluoranthene	22.86	252	10069	0.39	ng/ul	97
22) Benzo(k)fluoranthene	22.90	252	10849	0.39	ng/ul	98
23) Benzo(a)pyrene	23.43	252	9618	0.39	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.78	276	10146	0.39	ng/ul	99
25) Dibenzo(a,h)anthracene	25.80	278	8084	0.40	ng/ul	99
26) Benzo(g,h,i)perylene	26.47	276	8880	0.40	ng/ul	99

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

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