

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021115\
 Data File : BM000479.D
 Acq On : 11 Feb 2015 13:31
 Operator : TP/IZ
 Sample : SSTD0.801
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.801

Quant Time: Feb 11 14:13:01 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 13:45:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	2987	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	10231	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6342	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13911	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	12641	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	10360	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.10	152	10857	0.76	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	24304	0.76	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.55	128	19332	0.68	ng/ul	99
5) 2-Methylnaphthalene	12.17	142	13698	0.71	ng/ul	99
7) Acenaphthylene	14.09	152	20382	0.79	ng/ul	99
8) Acenaphthene	14.41	153	15589	0.77	ng/ul	99
9) Fluorene	15.42	166	18778	0.77	ng/ul	99
11) Pentachlorophenol	16.76	266	1129	1.11	ng/ul	100
12) Phenanthrene	17.15	178	30461	0.69	ng/ul	99
13) Anthracene	17.24	178	27799	0.70	ng/ul	99
15) Fluoranthene	19.18	202	35486	0.71	ng/ul	99
17) Pyrene	19.54	202	35558	0.65	ng/ul	100
18) Benzo(a)anthracene	21.28	228	28561	0.69	ng/ul	97
19) Chrysene	21.33	228	33106	0.68	ng/ul	99
21) Benzo(b)fluoranthene	22.86	252	28536	0.71	ng/ul	97
22) Benzo(k)fluoranthene	22.90	252	29589	0.70	ng/ul	96
23) Benzo(a)pyrene	23.43	252	26138	0.69	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.78	276	26820	0.63	ng/ul	99
25) Dibenzo(a,h)anthracene	25.79	278	21334	0.63	ng/ul	96
26) Benzo(g,h,i)perylene	26.47	276	23013	0.61	ng/ul	97

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

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