

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051115\
 Data File : BM001241.D
 Acq On : 11 May 2015 11:41
 Operator : TP/IZ
 Sample : SSTD0.459
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.459

Quant Time: May 12 02:41:48 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM051115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 12 02:23:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	1011	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	3406	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	1749	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	3656	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	2988	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	2686	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.84	152	1724	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	3168	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.29	128	3499	0.40	ng/ul	100
5) 2-Methylnaphthalene	11.91	142	2233	0.39	ng/ul	100
7) Acenaphthylene	13.84	152	3063	0.38	ng/ul	100
8) Acenaphthene	14.18	153	2430	0.39	ng/ul	100
9) Fluorene	15.19	166	2712	0.39	ng/ul	100
11) Pentachlorophenol	16.55	266	289	0.41	ng/ul	100
12) Phenanthrene	16.92	178	4109	0.39	ng/ul	100
13) Anthracene	17.01	178	3757	0.39	ng/ul	100
15) Fluoranthene	18.95	202	4612	0.38	ng/ul	100
17) Pyrene	19.31	202	4685	0.42	ng/ul	100
18) Benzo(a)anthracene	21.07	228	3602	0.38	ng/ul	100
19) Chrysene	21.12	228	4310	0.41	ng/ul	100
21) Benzo(b)fluoranthene	22.57	252	3869	0.41	ng/ul	100
22) Benzo(k)fluoranthene	22.61	252	4067	0.40	ng/ul	100
23) Benzo(a)pyrene	23.10	252	3656	0.39	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.29	276	4156	0.39	ng/ul	100
25) Dibenzo(a,h)anthracene	25.31	278	3326	0.39	ng/ul	100
26) Benzo(g,h,i)perylene	25.92	276	3746	0.40	ng/ul	100

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

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