

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030915\
 Data File : BM000664.D
 Acq On : 09 Mar 2015 17:10
 Operator : TP/IZ
 Sample : SSTD0.449
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.449

Quant Time: Mar 10 02:36:41 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM030915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 02:25:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	4033	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	13624	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	4117	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	12889	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	11165	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	10291	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	6421	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	10377	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	13564	0.39	ng/ul	100
5) 2-Methylnaphthalene	12.11	142	8574	0.35	ng/ul	100
7) Acenaphthylene	14.03	152	10597	0.55	ng/ul	100
8) Acenaphthene	14.36	153	8153	0.55	ng/ul	100
9) Fluorene	15.36	166	9581	0.53	ng/ul	100
11) Pentachlorophenol	16.71	266	295	0.09	ng/ul	100
12) Phenanthrene	17.10	178	14285	0.39	ng/ul	100
13) Anthracene	17.19	178	12682	0.37	ng/ul	100
15) Fluoranthene	19.11	202	15677	0.37	ng/ul	100
17) Pyrene	19.48	202	16374	0.33	ng/ul	100
18) Benzo(a)anthracene	21.23	228	12555	0.36	ng/ul	100
19) Chrysene	21.28	228	15445	0.40	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	14069	0.37	ng/ul	100
22) Benzo(k)fluoranthene	22.83	252	14932	0.39	ng/ul	100
23) Benzo(a)pyrene	23.35	252	13571	0.39	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.67	276	15592	0.45	ng/ul	100
25) Dibenzo(a,h)anthracene	25.68	278	12607	0.46	ng/ul	100
26) Benzo(g,h,i)perylene	26.35	276	13978	0.47	ng/ul	100

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

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