

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031017\
 Data File : BM009179.D
 Acq On : 10 Mar 2017 17:34
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.455

Quant Time: Mar 10 19:04:43 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 10 15:19:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	243	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	704	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	362	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	802	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	867	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	848	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	444	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	968	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.73	128	822	0.40	ng/ul#	89
5) 2-Methylnaphthalene	12.34	142	553	0.38	ng/ul	99
7) Acenaphthylene	14.23	152	778	0.41	ng/ul#	83
8) Acenaphthene	14.57	153	576	0.40	ng/ul	89
9) Fluorene	15.56	166	645	0.38	ng/ul#	80
11) Pentachlorophenol	16.92	266	148	0.38	ng/ul	98
12) Phenanthrene	17.30	178	998	0.40	ng/ul#	89
13) Anthracene	17.40	178	983	0.40	ng/ul#	95
15) Fluoranthene	19.32	202	1245	0.38	ng/ul	97
17) Pyrene	19.68	202	1246	0.42	ng/ul#	94
18) Benzo(a)anthracene	21.42	228	1221	0.42	ng/ul	97
19) Chrysene	21.47	228	1094	0.40	ng/ul	94
21) Benzo(b)fluoranthene	23.07	252	1332	0.40	ng/ul	91
22) Benzo(k)fluoranthene	23.11	252	1130	0.38	ng/ul	92
23) Benzo(a)pyrene	23.67	252	1210	0.40	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.15	276	1442	0.38	ng/ul#	90
25) Dibenzo(a,h)anthracene	26.18	278	1176	0.37	ng/ul	92
26) Benzo(g,h,i)perylene	26.89	276	1225	0.38	ng/ul#	90

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

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