

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092216\
 Data File : BM007520.D
 Acq On : 22 Sep 2016 20:28
 Operator : UM/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.442

Quant Time: Sep 23 12:15:13 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 23 04:00:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	1587	0.40	ng/ul	0.00
2) Naphthalene-d8	10.58	136	5967	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	3593	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.16	188	8939	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	9756	0.40	ng/ul	0.00
20) Perylene-d12	23.62	264	9182	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.17	152	2768	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.19	212	8344	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.62	128	5572	0.40	ng/ul	97
5) 2-Methylnaphthalene	12.25	142	3758	0.40	ng/ul	100
7) Acenaphthylene	14.14	152	5887	0.39	ng/ul	99
8) Acenaphthene	14.49	153	4679	0.39	ng/ul	97
9) Fluorene	15.48	166	5679	0.39	ng/ul	95
11) Pentachlorophenol	16.83	266	963	0.39	ng/ul	97
12) Phenanthrene	17.21	178	9741	0.40	ng/ul	98
13) Anthracene	17.30	178	8744	0.40	ng/ul	98
15) Fluoranthene	19.22	202	12306	0.40	ng/ul	91
17) Pyrene	19.59	202	13403	0.42	ng/ul	99
18) Benzo(a)anthracene	21.32	228	11893	0.41	ng/ul	93
19) Chrysene	21.38	228	12444	0.40	ng/ul	99
21) Benzo(b)fluoranthene	22.93	252	12944	0.42	ng/ul	99
22) Benzo(k)fluoranthene	22.97	252	12302	0.40	ng/ul	99
23) Benzo(a)pyrene	23.51	252	11848	0.42	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.91	276	12566	0.43	ng/ul	100
25) Dibenzo(a,h)anthracene	25.92	278	9903	0.44	ng/ul	100
26) Benzo(g,h,i)perylene	26.61	276	11129	0.42	ng/ul	99

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

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