

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093223.D
 Acq On : 16 Jun 2017 00:12
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.409

Quant Time: Jun 16 04:45:56 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 04:43:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2379	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	12018	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6621	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	15697	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16963	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	14282	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7509	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	17908	0.42	ng/ul	0.00

Target Compounds					Qvalue	
3) Naphthalene	10.57	128	12149	0.407	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	8686	0.426	ng/ul	99
7) Acenaphthylene	14.08	152	12757	0.449	ng/ul	96
8) Acenaphthene	14.43	153	9486	0.414	ng/ul	99
9) Fluorene	15.42	166	10512	0.391	ng/ul#	90
11) Pentachlorophenol	16.75	266	1035	0.444	ng/ul	95
12) Phenanthrene	17.13	178	18159	0.420	ng/ul#	87
13) Anthracene	17.22	178	17440	0.464	ng/ul#	87
15) Fluoranthene	19.16	202	21953	0.414	ng/ul	83
17) Pyrene	19.52	202	23488	0.470	ng/ul#	83
18) Benzo(a)anthracene	21.25	228	20645	0.456	ng/ul#	86
19) Chrysene	21.30	228	19661	0.419	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	18643	0.430	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	17655	0.384	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	17421	0.425	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.18	276	19779	0.417	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.21	278	16456	0.412	ng/ul#	89
26) Benzo(g,h,i)perylene	26.96	276	16284	0.400	ng/ul	97

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

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