

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061917\
 Data File : BE093243.D
 Acq On : 19 Jun 2017 10:27
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.413

Quant Time: Jun 19 10:59:24 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 19 07:31:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2206	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	11624	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6958	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	18212	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	19701	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	16906	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7495	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	20935	0.42	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.57	128	11814	0.410	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	8629	0.438	ng/ul	98
7) Acenaphthylene	14.09	152	13285	0.445	ng/ul	96
8) Acenaphthene	14.43	153	9910	0.411	ng/ul	97
9) Fluorene	15.42	166	11347	0.402	ng/ul#	93
11) Pentachlorophenol	16.77	266	1132	0.419	ng/ul	96
12) Phenanthrene	17.13	178	20141	0.402	ng/ul#	87
13) Anthracene	17.22	178	19449	0.446	ng/ul#	87
15) Fluoranthene	19.15	202	24919	0.406	ng/ul	83
17) Pyrene	19.51	202	26657	0.459	ng/ul#	86
18) Benzo(a)anthracene	21.25	228	23997	0.457	ng/ul#	86
19) Chrysene	21.30	228	22767	0.418	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	22014	0.429	ng/ul	90
22) Benzo(k)fluoranthene	22.98	252	21044	0.386	ng/ul#	91
23) Benzo(a)pyrene	23.55	252	20534	0.423	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.19	276	23370	0.416	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.21	278	19286	0.408	ng/ul	90
26) Benzo(g,h,i)perylene	26.96	276	19424	0.403	ng/ul	96

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

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