

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061317\
 Data File : BE093144.D
 Acq On : 13 Jun 2017 23:22
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.402

Quant Time: Jun 14 01:27:00 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 14 01:25:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3600	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	16492	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	9290	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	24485	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	29137	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	24223	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	10146	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	28056	0.42	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.57	128	16877	0.412	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	11719	0.419	ng/ul	99
7) Acenaphthylene	14.08	152	16241	0.407	ng/ul	96
8) Acenaphthene	14.43	153	13128	0.408	ng/ul	99
9) Fluorene	15.42	166	15090	0.400	ng/ul#	91
11) Pentachlorophenol	16.75	266	1779	0.490	ng/ul	97
12) Phenanthrene	17.14	178	28455	0.422	ng/ul#	87
13) Anthracene	17.22	178	25795	0.440	ng/ul#	88
15) Fluoranthene	19.15	202	34260	0.415	ng/ul	84
17) Pyrene	19.51	202	35963	0.419	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	34018	0.438	ng/ul#	85
19) Chrysene	21.30	228	33414	0.414	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	30720	0.418	ng/ul	88
22) Benzo(k)fluoranthene	22.97	252	31370	0.402	ng/ul#	89
23) Benzo(a)pyrene	23.55	252	29522	0.425	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.18	276	32944	0.409	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.20	278	27300	0.403	ng/ul#	91
26) Benzo(g,h,i)perylene	26.95	276	27617	0.400	ng/ul	97

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

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