

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061917\
 Data File : BE093274.D
 Acq On : 20 Jun 2017 5:39
 Operator : SJ/MA
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.415

Quant Time: Jun 20 08:51:58 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 08:49:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1895	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	10225	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5841	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14645	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	15175	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	13151	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	6476	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	16913	0.42	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	10376	0.409	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	7470	0.431	ng/ul	99
7) Acenaphthylene	14.09	152	12356	0.493	ng/ul	97
8) Acenaphthene	14.42	153	8355	0.413	ng/ul	97
9) Fluorene	15.42	166	9750	0.411	ng/ul#	94
11) Pentachlorophenol	16.75	266	1259	0.579	ng/ul	99
12) Phenanthrene	17.14	178	16552	0.411	ng/ul#	88
13) Anthracene	17.22	178	16978	0.484	ng/ul#	89
15) Fluoranthene	19.15	202	20019	0.405	ng/ul	84
17) Pyrene	19.51	202	21183	0.473	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	18720	0.462	ng/ul#	83
19) Chrysene	21.30	228	17146	0.408	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	16925	0.424	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	16045	0.379	ng/ul	91
23) Benzo(a)pyrene	23.55	252	16127	0.427	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.18	276	18015	0.412	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.20	278	15023	0.408	ng/ul#	87
26) Benzo(g,h,i)perylene	26.96	276	14876	0.397	ng/ul#	93

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

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