

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080417\
 Data File : BE093670.D
 Acq On : 4 Aug 2017 14:59
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.443

Quant Time: Aug 04 17:05:03 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 04 16:09:39 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2607	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	13428	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7726	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	18579	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	19942	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	18863	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8570	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	21258	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	13150	0.403	ng/ul#	91
5) 2-Methylnaphthalene	12.36	142	9273	0.389	ng/ul	99
7) Acenaphthylene	14.25	152	13647	0.422	ng/ul	96
8) Acenaphthene	14.59	153	10165	0.400	ng/ul	96
9) Fluorene	15.57	166	11999	0.373	ng/ul#	95
11) Pentachlorophenol	16.91	266	1090	0.356	ng/ul	96
12) Phenanthrene	17.29	178	19877	0.393	ng/ul#	88
13) Anthracene	17.38	178	19499	0.396	ng/ul#	88
15) Fluoranthene	19.30	202	22403	0.354	ng/ul	83
17) Pyrene	19.66	202	23253	0.399	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	22924	0.396	ng/ul#	86
19) Chrysene	21.45	228	21782	0.393	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	21984	0.377	ng/ul	87
22) Benzo(k)fluoranthene	23.20	252	21663	0.376	ng/ul#	88
23) Benzo(a)pyrene	23.80	252	21334	0.384	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.55	276	23357	0.381	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.57	278	19528	0.381	ng/ul#	91
26) Benzo(g,h,i)perylene	27.37	276	19226	0.372	ng/ul	99

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

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