

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080817\
 Data File : BE093709.D
 Acq On : 8 Aug 2017 10:02
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.444

Quant Time: Aug 09 04:13:35 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	7.92	152	2771	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	13770	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7904	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	19349	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	21299	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19789	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8758	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	21714	0.36	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	13524	0.404	ng/ul#	91
5) 2-Methylnaphthalene	12.36	142	9502	0.389	ng/ul	98
7) Acenaphthylene	14.24	152	13113	0.397	ng/ul#	95
8) Acenaphthene	14.59	153	10279	0.395	ng/ul	99
9) Fluorene	15.58	166	12124	0.369	ng/ul#	95
11) Pentachlorophenol	16.93	266	860	0.269	ng/ul	98
12) Phenanthrene	17.29	178	20973	0.398	ng/ul#	87
13) Anthracene	17.38	178	19628	0.382	ng/ul#	87
15) Fluoranthene	19.30	202	23488	0.356	ng/ul	84
17) Pyrene	19.66	202	24244	0.390	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	23221	0.376	ng/ul#	86
19) Chrysene	21.45	228	23689	0.400	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	23243	0.380	ng/ul	85
22) Benzo(k)fluoranthene	23.19	252	22905	0.379	ng/ul#	87
23) Benzo(a)pyrene	23.80	252	22213	0.381	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.55	276	24665	0.384	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.57	278	20498	0.382	ng/ul#	89
26) Benzo(g,h,i)perylene	27.37	276	20141	0.372	ng/ul	96

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

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