

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080417\
 Data File : BE093662.D
 Acq On : 4 Aug 2017 8:49
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.442

Quant Time: Aug 04 16:09:24 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	2342	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	11609	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	6606	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	16803	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	20891	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19010	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7259	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	20572	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	11590	0.410	ng/ul#	91
5) 2-Methylnaphthalene	12.36	142	7949	0.386	ng/ul	98
7) Acenaphthylene	14.25	152	11015	0.399	ng/ul	96
8) Acenaphthene	14.59	153	8660	0.398	ng/ul	98
9) Fluorene	15.58	166	10249	0.373	ng/ul#	95
11) Pentachlorophenol	16.93	266	781	0.282	ng/ul	95
12) Phenanthrene	17.29	178	18323	0.400	ng/ul#	88
13) Anthracene	17.38	178	16904	0.379	ng/ul#	88
15) Fluoranthene	19.30	202	22080	0.386	ng/ul	84
17) Pyrene	19.66	202	23072	0.378	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	23505	0.388	ng/ul#	85
19) Chrysene	21.45	228	22988	0.396	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	22665	0.385	ng/ul	85
22) Benzo(k)fluoranthene	23.19	252	22024	0.380	ng/ul#	85
23) Benzo(a)pyrene	23.80	252	21646	0.387	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.55	276	23804	0.385	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.57	278	19737	0.382	ng/ul#	89
26) Benzo(g,h,i)perylene	27.36	276	19855	0.382	ng/ul	96

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

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