

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081017\
 Data File : BE093740.D
 Acq On : 10 Aug 2017 9:39
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.447

Quant Time: Aug 11 17:29:14 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 11 06:07:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2236	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	11713	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	6955	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	17133	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	18614	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	17003	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7600	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	19203	0.36	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.75	128	11542	0.405	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	8200	0.395	ng/ul	98
7) Acenaphthylene	14.24	152	11640	0.400	ng/ul	97
8) Acenaphthene	14.59	153	9123	0.399	ng/ul	98
9) Fluorene	15.57	166	10788	0.373	ng/ul#	90
11) Pentachlorophenol	16.93	266	837	0.296	ng/ul	96
12) Phenanthrene	17.29	178	18513	0.397	ng/ul#	88
13) Anthracene	17.38	178	17285	0.380	ng/ul#	88
15) Fluoranthene	19.30	202	20916	0.358	ng/ul	84
17) Pyrene	19.66	202	21527	0.396	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	20538	0.381	ng/ul#	85
19) Chrysene	21.44	228	20615	0.398	ng/ul	97
21) Benzo(b)fluoranthene	23.14	252	20633	0.392	ng/ul	86
22) Benzo(k)fluoranthene	23.19	252	19541	0.377	ng/ul#	86
23) Benzo(a)pyrene	23.80	252	19612	0.392	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.55	276	21109	0.382	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.57	278	17489	0.379	ng/ul#	91
26) Benzo(g,h,i)perylene	27.36	276	17235	0.370	ng/ul	97

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

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