

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093592.D
 Acq On : 26 Jul 2017 00:11
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.440

Quant Time: Jul 26 02:50:16 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 06:10:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2144	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10734	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	6670	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	16745	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	18955	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	17392	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7045	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	19733	0.37	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.76	128	10617	0.407	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	7683	0.403	ng/ul	98
7) Acenaphthylene	14.25	152	11362	0.407	ng/ul	96
8) Acenaphthene	14.59	153	8788	0.400	ng/ul	98
9) Fluorene	15.57	166	10580	0.381	ng/ul#	94
11) Pentachlorophenol	16.91	266	786	0.285	ng/ul	96
12) Phenanthrene	17.29	178	18713	0.410	ng/ul#	88
13) Anthracene	17.38	178	17344	0.390	ng/ul#	88
15) Fluoranthene	19.30	202	21208	0.372	ng/ul	85
17) Pyrene	19.66	202	21831	0.395	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	21173	0.385	ng/ul#	86
19) Chrysene	21.45	228	20753	0.394	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	20823	0.387	ng/ul	86
22) Benzo(k)fluoranthene	23.19	252	20106	0.379	ng/ul#	87
23) Benzo(a)pyrene	23.80	252	19834	0.388	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.55	276	21814	0.386	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.57	278	18243	0.386	ng/ul#	90
26) Benzo(g,h,i)perylene	27.35	276	18061	0.379	ng/ul	96

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

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