

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093557.D
 Acq On : 25 Jul 2017 2:39
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.437

Quant Time: Jul 25 05:45:19 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 24 18:07:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2165	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	11257	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7354	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	20468	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	23605	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	21379	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7348	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	24475	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	11120	0.406	ng/ul#	92
5) 2-Methylnaphthalene	12.37	142	8119	0.406	ng/ul	97
7) Acenaphthylene	14.25	152	12226	0.397	ng/ul	96
8) Acenaphthene	14.59	153	9614	0.397	ng/ul	97
9) Fluorene	15.57	166	12063	0.394	ng/ul#	92
11) Pentachlorophenol	16.93	266	1045	0.310	ng/ul	95
12) Phenanthrene	17.29	178	21714	0.390	ng/ul#	86
13) Anthracene	17.39	178	20737	0.382	ng/ul#	92
15) Fluoranthene	19.31	202	26396	0.379	ng/ul	84
17) Pyrene	19.67	202	26993	0.392	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	26290	0.384	ng/ul#	85
19) Chrysene	21.45	228	25896	0.394	ng/ul	97
21) Benzo(b)fluoranthene	23.14	252	25582	0.387	ng/ul	85
22) Benzo(k)fluoranthene	23.19	252	24749	0.379	ng/ul#	88
23) Benzo(a)pyrene	23.80	252	24297	0.386	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.55	276	26790	0.386	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.57	278	22602	0.389	ng/ul#	90
26) Benzo(g,h,i)perylene	27.37	276	22166	0.379	ng/ul	97

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

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