

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093540.D
 Acq On : 24 Jul 2017 16:13
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.436

Quant Time: Jul 24 17:38:07 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 13:14:05 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2754	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	13789	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	8965	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	24494	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	28008	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	24935	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	9101	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	29273	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	13525	0.403	ng/ul#	91
5) 2-Methylnaphthalene	12.37	142	9809	0.401	ng/ul	99
7) Acenaphthylene	14.25	152	14934	0.398	ng/ul	96
8) Acenaphthene	14.59	153	11784	0.399	ng/ul	96
9) Fluorene	15.58	166	14742	0.395	ng/ul#	92
11) Pentachlorophenol	16.93	266	1245	0.308	ng/ul	97
12) Phenanthrene	17.29	178	26467	0.397	ng/ul#	85
13) Anthracene	17.39	178	24893	0.383	ng/ul#	92
15) Fluoranthene	19.31	202	31238	0.374	ng/ul	84
17) Pyrene	19.67	202	31920	0.390	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	31063	0.383	ng/ul#	86
19) Chrysene	21.45	228	30668	0.394	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	30195	0.391	ng/ul	86
22) Benzo(k)fluoranthene	23.20	252	29332	0.386	ng/ul#	88
23) Benzo(a)pyrene	23.81	252	28336	0.386	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.56	276	31430	0.388	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.58	278	26537	0.392	ng/ul#	90
26) Benzo(g,h,i)perylene	27.38	276	26339	0.386	ng/ul	97

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

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