

Data Path : Z:\HPCHEM1\BNA_E\Data\BE072417\
 Data File : BE093574.D
 Acq On : 25 Jul 2017 13:06
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.438

Quant Time: Jul 25 13:46:50 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	7.93	152	1890	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	9678	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	6248	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	17502	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	20321	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	18645	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6412	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	21003	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	9546	0.405	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	6957	0.405	ng/ul	98
7) Acenaphthylene	14.25	152	10549	0.404	ng/ul	96
8) Acenaphthene	14.59	153	8269	0.402	ng/ul	99
9) Fluorene	15.58	166	10438	0.402	ng/ul#	93
11) Pentachlorophenol	16.93	266	864	0.299	ng/ul	95
12) Phenanthrene	17.29	178	18756	0.393	ng/ul#	86
13) Anthracene	17.38	178	17635	0.380	ng/ul#	86
15) Fluoranthene	19.30	202	22444	0.377	ng/ul	84
17) Pyrene	19.67	202	23221	0.391	ng/ul#	79
18) Benzo(a)anthracene	21.39	228	22628	0.384	ng/ul#	86
19) Chrysene	21.45	228	22317	0.395	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	21858	0.379	ng/ul	86
22) Benzo(k)fluoranthene	23.19	252	21869	0.385	ng/ul#	86
23) Benzo(a)pyrene	23.80	252	21124	0.385	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.55	276	23318	0.385	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.57	278	19591	0.387	ng/ul#	91
26) Benzo(g,h,i)perylene	27.37	276	19386	0.380	ng/ul	96

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

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