

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081517\
 Data File : BE093831.D
 Acq On : 16 Aug 2017 5:44
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.408

Quant Time: Aug 16 07:39:24 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 07:30:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1920	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10201	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5790	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13246	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	14742	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	12663	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6371	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	15694	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	9807	0.395	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	6770	0.374	ng/ul	99
7) Acenaphthylene	14.24	152	10103	0.417	ng/ul	97
8) Acenaphthene	14.59	153	7418	0.389	ng/ul	98
9) Fluorene	15.57	166	8561	0.355	ng/ul#	89
11) Pentachlorophenol	16.93	266	929	0.425	ng/ul	97
12) Phenanthrene	17.29	178	13883	0.385	ng/ul#	90
13) Anthracene	17.38	178	13912	0.396	ng/ul#	89
15) Fluoranthene	19.30	202	16404	0.364	ng/ul	85
17) Pyrene	19.66	202	17430	0.405	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	16556	0.387	ng/ul#	87
19) Chrysene	21.45	228	16406	0.400	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	14416	0.368	ng/ul	88
22) Benzo(k)fluoranthene	23.20	252	15622	0.404	ng/ul#	90
23) Benzo(a)pyrene	23.80	252	14701	0.395	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.56	276	14237	0.346	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.58	278	11869	0.345	ng/ul	95
26) Benzo(g,h,i)perylene	27.37	276	11624	0.335	ng/ul	99

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

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