

Data Path : Z:\HPCHEM1\BNA_E\Data\BE081417\
 Data File : BE093783.D
 Acq On : 14 Aug 2017 10:20
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.403

Quant Time: Aug 14 18:13:47 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 12 02:01:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2117	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10420	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5896	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14300	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	16344	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	14501	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6425	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	16393	0.36	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	10116	0.399	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	6947	0.376	ng/ul	98
7) Acenaphthylene	14.24	152	9477	0.384	ng/ul	97
8) Acenaphthene	14.59	153	7436	0.383	ng/ul	97
9) Fluorene	15.57	166	8816	0.359	ng/ul#	91
11) Pentachlorophenol	16.93	266	648	0.275	ng/ul	97
12) Phenanthrene	17.29	178	15449	0.397	ng/ul#	89
13) Anthracene	17.38	178	14429	0.380	ng/ul#	90
15) Fluoranthene	19.30	202	17840	0.366	ng/ul	85
17) Pyrene	19.66	202	18413	0.386	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	17628	0.372	ng/ul#	87
19) Chrysene	21.45	228	17992	0.396	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	17101	0.381	ng/ul	85
22) Benzo(k)fluoranthene	23.19	252	17325	0.392	ng/ul#	87
23) Benzo(a)pyrene	23.80	252	16363	0.383	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.56	276	17210	0.365	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.58	278	14109	0.358	ng/ul#	90
26) Benzo(g,h,i)perylene	27.37	276	14573	0.367	ng/ul	99

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

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