

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098093.D
 Acq On : 26 Oct 2018 17:53
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.472

Quant Time: Oct 26 23:21:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 26 23:15:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	915	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4628	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3300	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8062	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	12161	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	10109	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3165	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	9729	0.37	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.73	128	4301	0.388	ng/ul	99
5) 2-Methylnaphthalene	12.34	142	3237	0.387	ng/ul	100
7) Acenaphthylene	14.25	152	4939	0.379	ng/ul	99
8) Acenaphthene	14.60	153	3658	0.376	ng/ul	98
9) Fluorene	15.59	166	4381	0.374	ng/ul	99
11) Pentachlorophenol	16.95	266	208	0.493	ng/ul	97
12) Phenanthrene	17.33	178	7060	0.379	ng/ul	100
13) Anthracene	17.42	178	6938	0.378	ng/ul	98
15) Fluoranthene	19.35	202	8814	0.372	ng/ul	99
17) Pyrene	19.71	202	9027	0.318	ng/ul	100
18) Benzo(a)anthracene	21.45	228	11428	0.320	ng/ul	99
19) Chrysene	21.51	228	9579	0.346	ng/ul	98
21) Benzo(b)fluoranthene	23.24	252	9598	0.343	ng/ul	99
22) Benzo(k)fluoranthene	23.29	252	9730	0.355	ng/ul	99
23) Benzo(a)pyrene	23.90	252	9387	0.348	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.71	276	10933	0.354	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.74	278	8990	0.349	ng/ul	97
26) Benzo(g,h,i)perylene	27.54	276	9055	0.348	ng/ul	99

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

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