

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102218\
 Data File : BE098013.D
 Acq On : 22 Oct 2018 17:28
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.459

Quant Time: Oct 23 01:28:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 01:24:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	1567	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	7541	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	5261	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	13166	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	17116	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	14777	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	5111	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	15929	0.37	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.73	128	6996	0.385	ng/ul	99
5) 2-Methylnaphthalene	12.35	142	5222	0.433	ng/ul	97
7) Acenaphthylene	14.26	152	8176	0.441	ng/ul	98
8) Acenaphthene	14.60	153	6004	0.389	ng/ul	97
9) Fluorene	15.59	166	7211	0.387	ng/ul	100
11) Pentachlorophenol	16.96	266	656	0.546	ng/ul	96
12) Phenanthrene	17.34	178	11782	0.375	ng/ul	100
13) Anthracene	17.43	178	11716	0.439	ng/ul	100
15) Fluoranthene	19.36	202	14571	0.351	ng/ul	99
17) Pyrene	19.72	202	16998	0.404	ng/ul	99
18) Benzo(a)anthracene	21.46	228	19206	0.404	ng/ul	99
19) Chrysene	21.51	228	15447	0.337	ng/ul	97
21) Benzo(b)fluoranthene	23.24	252	15548	0.407	ng/ul	97
22) Benzo(k)fluoranthene	23.30	252	15047	0.348	ng/ul	95
23) Benzo(a)pyrene	23.91	252	14810	0.424	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	26.72	276	17256	0.435	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.75	278	14260	0.430	ng/ul	96
26) Benzo(g,h,i)perylene	27.55	276	14528	0.388	ng/ul	97

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

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