

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098113.D
 Acq On : 31 Oct 2018 11:12
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.476

Quant Time: Oct 31 11:56:49 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 : Tue Oct 30 01:14:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	1023	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5042	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	3487	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8348	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	9619	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9514	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3529	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10593	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.73	128	4815	0.399	ng/ul	100
5) 2-Methylnaphthalene	12.35	142	3663	0.401	ng/ul	99
7) Acenaphthylene	14.26	152	5429	0.395	ng/ul	98
8) Acenaphthene	14.60	153	4072	0.396	ng/ul	98
9) Fluorene	15.59	166	4672	0.378	ng/ul	100
11) Pentachlorophenol	16.95	266	215	0.492	ng/ul	94
12) Phenanthrene	17.33	178	7706	0.400	ng/ul	99
13) Anthracene	17.42	178	7425	0.390	ng/ul	99
15) Fluoranthene	19.35	202	9619	0.392	ng/ul	97
17) Pyrene	19.71	202	9719	0.433	ng/ul	98
18) Benzo(a)anthracene	21.45	228	12311	0.435	ng/ul	99
19) Chrysene	21.51	228	10484	0.479	ng/ul	100
21) Benzo(b)fluoranthene	23.24	252	10839	0.411	ng/ul	99
22) Benzo(k)fluoranthene	23.29	252	10128	0.393	ng/ul	99
23) Benzo(a)pyrene	23.91	252	10033	0.396	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.72	276	11375	0.392	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.75	278	9475	0.391	ng/ul#	96
26) Benzo(g,h,i)perylene	27.54	276	9636	0.394	ng/ul	98

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

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