

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098053.D
 Acq On : 25 Oct 2018 13:02
 Operator : SJ/JU
 Sample : SST00.864
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.864

Quant Time: Oct 25 13:31:49 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 : Thu Oct 25 13:24:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	929	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4761	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3279	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8079	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11726	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9559	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	6986	0.94	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	21706	0.82	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.73	128	9269	0.809	ng/ul	98
5) 2-Methylnaphthalene	12.34	142	7052	0.926	ng/ul	100
7) Acenaphthylene	14.25	152	10899	0.943	ng/ul	98
8) Acenaphthene	14.60	153	7934	0.824	ng/ul	98
9) Fluorene	15.59	166	9767	0.842	ng/ul	99
11) Pentachlorophenol	16.95	266	320	0.434	ng/ul	96
12) Phenanthrene	17.33	178	15370	0.796	ng/ul	99
13) Anthracene	17.42	178	15329	0.935	ng/ul	100
15) Fluoranthene	19.35	202	19674	0.773	ng/ul	99
17) Pyrene	19.71	202	23152	0.804	ng/ul	98
18) Benzo(a)anthracene	21.45	228	26508	0.814	ng/ul	99
19) Chrysene	21.51	228	21415	0.682	ng/ul	99
21) Benzo(b)fluoranthene	23.24	252	21044	0.852	ng/ul	97
22) Benzo(k)fluoranthene	23.29	252	21995	0.786	ng/ul	97
23) Benzo(a)pyrene	23.91	252	20793	0.920	ng/ul#	98
24) Indeno(1,2,3-cd)pyrene	26.71	276	24195	0.943	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.73	278	20079	0.936	ng/ul	97
26) Benzo(g,h,i)perylene	27.54	276	20386	0.842	ng/ul	98

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

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