

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098051.D
 Acq On : 25 Oct 2018 11:47
 Operator : SJ/JU
 Sample : SST0.262
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.262

Quant Time: Oct 25 13:29:53 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	7.87	152	918	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4637	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	8248	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11685	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9578	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1631	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	5083	0.19	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	2148	0.192	ng/ul	94
5) 2-Methylnaphthalene	12.34	142	1641	0.221	ng/ul	99
7) Acenaphthylene	14.25	152	2508	0.216	ng/ul	98
8) Acenaphthene	14.60	153	1905	0.197	ng/ul	98
9) Fluorene	15.59	166	2272	0.195	ng/ul	99
11) Pentachlorophenol	16.95	266	66	0.088	ng/ul#	82
12) Phenanthrene	17.33	178	3600	0.183	ng/ul	96
13) Anthracene	17.42	178	3593	0.215	ng/ul	97
15) Fluoranthene	19.35	202	4635	0.178	ng/ul	97
17) Pyrene	19.71	202	4702	0.164	ng/ul	98
18) Benzo(a)anthracene	21.45	228	6890	0.212	ng/ul	96
19) Chrysene	21.51	228	5084	0.162	ng/ul	99
21) Benzo(b)fluoranthene	23.24	252	5296	0.214	ng/ul	92
22) Benzo(k)fluoranthene	23.29	252	4962	0.177	ng/ul#	95
23) Benzo(a)pyrene	23.91	252	4935	0.218	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	26.71	276	5700	0.222	ng/ul#	64
25) Dibenzo(a,h)anthracene	26.74	278	4663	0.217	ng/ul#	88
26) Benzo(g,h,i)perylene	27.54	276	4761	0.196	ng/ul	94

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

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