

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097850.D
 Acq On : 11 Oct 2018 10:21
 Operator : SJ/JU
 Sample : SST00.144
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.144

Quant Time: Oct 11 12:21:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 12:14:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2202	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10360	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	7850	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	24068	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	33712	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	28679	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1713	0.11	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	7433	0.09	ng/ul	0.00

Target Compounds					Qvalue	
3) Naphthalene	10.74	128	2471	0.099	ng/ul#	87
5) 2-Methylnaphthalene	12.36	142	1792	0.108	ng/ul	99
7) Acenaphthylene	14.26	152	2853	0.103	ng/ul	97
8) Acenaphthene	14.60	153	2268	0.098	ng/ul	98
9) Fluorene	15.60	166	2903	0.105	ng/ul	98
12) Phenanthrene	17.34	178	5491	0.095	ng/ul	96
13) Anthracene	17.43	178	4995	0.102	ng/ul	97
15) Fluoranthene	19.36	202	6856	0.090	ng/ul	95
17) Pyrene	19.72	202	8722	0.105	ng/ul#	96
18) Benzo(a)anthracene	21.46	228	9599	0.103	ng/ul	99
19) Chrysene	21.52	228	8281	0.092	ng/ul	96
21) Benzo(b)fluoranthene	23.25	252	7335	0.099	ng/ul	89
22) Benzo(k)fluoranthene	23.30	252	8022	0.095	ng/ul#	85
23) Benzo(a)pyrene	23.91	252	7418	0.109	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.72	276	8284	0.108	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.75	278	6834	0.106	ng/ul#	86
26) Benzo(g,h,i)perylene	27.54	276	7352	0.101	ng/ul#	92

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

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