

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080118\
 Data File : BE097128.D
 Acq On : 1 Aug 2018 13:42
 Operator : SJ/JU
 Sample : SST0.475
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.475

Quant Time: Aug 01 14:22:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 01 14:22:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1463	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7065	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3864	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10474	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	13426	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	13532	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4539	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13933	0.53	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	6850	0.401	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	4502	0.387	ng/ul	97
7) Acenaphthylene	13.74	152	6676	0.422	ng/ul	97
8) Acenaphthene	14.08	153	4882	0.373	ng/ul	97
9) Fluorene	15.08	166	5558	0.412	ng/ul	93
11) Pentachlorophenol	16.44	266	536	0.782	ng/ul	93
12) Phenanthrene	16.81	178	10531	0.390	ng/ul#	89
13) Anthracene	16.90	178	9802	0.427	ng/ul#	92
15) Fluoranthene	18.84	202	13653	0.475	ng/ul	84
17) Pyrene	19.20	202	14276	0.324	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	15032	0.475	ng/ul#	84
19) Chrysene	21.01	228	15664	0.382	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	14405	0.390	ng/ul	85
22) Benzo(k)fluoranthene	22.58	252	15459	0.383	ng/ul#	88
23) Benzo(a)pyrene	23.11	252	14531	0.465	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.50	276	18541	0.572	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.52	278	15113	0.587	ng/ul#	93
26) Benzo(g,h,i)perylene	26.20	276	16126	0.512	ng/ul#	94

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

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