

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080118\
 Data File : BE097127.D
 Acq On : 1 Aug 2018 13:03
 Operator : SJ/JU
 Sample : SST0.274
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.274

Quant Time: Aug 01 14:22:24 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	7.39	152	1446	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6710	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3685	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9848	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12892	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	13474	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	2155	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	6810	0.28	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	10.18	128	3238	0.200	ng/ul#	96
5) 2-Methylnaphthalene	11.81	142	2120	0.192	ng/ul	98
7) Acenaphthylene	13.73	152	3158	0.209	ng/ul	99
8) Acenaphthene	14.08	153	2272	0.182	ng/ul#	91
9) Fluorene	15.08	166	2575	0.200	ng/ul	90
11) Pentachlorophenol	16.44	266	268	0.416	ng/ul	98
12) Phenanthrene	16.81	178	5018	0.198	ng/ul#	89
13) Anthracene	16.90	178	4582	0.213	ng/ul#	92
15) Fluoranthene	18.84	202	6580	0.243	ng/ul	84
17) Pyrene	19.20	202	6998	0.165	ng/ul#	82
18) Benzo(a)anthracene	20.96	228	7426	0.244	ng/ul#	86
19) Chrysene	21.01	228	7646	0.194	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	7185	0.195	ng/ul	90
22) Benzo(k)fluoranthene	22.58	252	7806	0.194	ng/ul	91
23) Benzo(a)pyrene	23.11	252	7349	0.236	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.50	276	9097	0.282	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.53	278	7356	0.287	ng/ul#	83
26) Benzo(g,h,i)perylene	26.21	276	7841	0.250	ng/ul	97

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

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