

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097291.D
 Acq On : 17 Aug 2018 11:01
 Operator : SJ/JU
 Sample : SST00.202
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.202

Quant Time: Aug 17 12:20:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 12:19:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	925	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	4677	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	2404	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	5449	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	4447	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	4032	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	1464	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	2806	0.15	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	2280	0.201	ng/ul#	95
5) 2-Methylnaphthalene	11.81	142	1426	0.191	ng/ul	99
7) Acenaphthylene	13.74	152	2197	0.211	ng/ul	99
8) Acenaphthene	14.08	153	1541	0.205	ng/ul#	89
9) Fluorene	15.08	166	1677	0.194	ng/ul	93
11) Pentachlorophenol	16.46	266	56	0.072	ng/ul#	81
12) Phenanthrene	16.81	178	2809	0.202	ng/ul#	89
13) Anthracene	16.90	178	2249	0.174	ng/ul#	91
15) Fluoranthene	18.84	202	2843	0.155	ng/ul	85
17) Pyrene	19.21	202	2827	0.237	ng/ul#	84
18) Benzo(a)anthracene	20.96	228	2108	0.166	ng/ul#	84
19) Chrysene	21.02	228	2711	0.207	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	1999	0.182	ng/ul	91
22) Benzo(k)fluoranthene	22.58	252	2671	0.225	ng/ul	93
23) Benzo(a)pyrene	23.12	252	1905	0.173	ng/ul	90
24) Indeno(1,2,3-cd)pyrene	25.53	276	2374	0.172	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.55	278	1878	0.167	ng/ul#	95
26) Benzo(g,h,i)perylene	26.22	276	1941	0.163	ng/ul	94

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

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