

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122818\
 Data File : BE098568.D
 Acq On : 28 Dec 2018 11:13
 Operator : SJ/JU
 Sample : SST0.805
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.805

Quant Time: Dec 28 11:45:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 11:34:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	838	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	3978	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	2532	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	6457	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	7833	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	7557	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.21	152	5385	0.76	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	16983	0.81	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	8837	0.928	ng/ul	97
5) 2-Methylnaphthalene	12.28	142	5874	0.816	ng/ul	100
7) Acenaphthylene	14.18	152	8987	0.900	ng/ul	98
8) Acenaphthene	14.53	153	7118	0.954	ng/ul	98
9) Fluorene	15.52	166	8562	0.953	ng/ul	99
11) Pentachlorophenol	16.89	266	709	2.099	ng/ul	95
12) Phenanthrene	17.27	178	14356	0.963	ng/ul	99
13) Anthracene	17.36	178	12558	0.853	ng/ul	99
15) Fluoranthene	19.29	202	18595	0.979	ng/ul	94
17) Pyrene	19.65	202	21148	1.156	ng/ul#	95
18) Benzo(a)anthracene	21.39	228	18777	0.815	ng/ul	99
19) Chrysene	21.45	228	22304	1.252	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	19135	0.914	ng/ul	97
22) Benzo(k)fluoranthene	23.21	252	20215	0.986	ng/ul	98
23) Benzo(a)pyrene	23.81	252	18173	0.902	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.57	276	21202	0.919	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.59	278	17399	0.905	ng/ul	98
26) Benzo(g,h,i)perylene	27.38	276	18297	0.942	ng/ul	96

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

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