

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122818\
 Data File : BE098567.D
 Acq On : 28 Dec 2018 10:37
 Operator : SJ/JU
 Sample : SST00.404
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.404

Quant Time: Dec 28 11:37:33 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	885	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	4236	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	2684	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	6759	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	7989	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	7966	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.21	152	2699	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	8387	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	4577	0.451	ng/ul	100
5) 2-Methylnaphthalene	12.28	142	2845	0.371	ng/ul	98
7) Acenaphthylene	14.19	152	4428	0.418	ng/ul	99
8) Acenaphthene	14.53	153	3537	0.447	ng/ul	98
9) Fluorene	15.53	166	4231	0.444	ng/ul	99
11) Pentachlorophenol	16.89	266	331	0.936	ng/ul	95
12) Phenanthrene	17.27	178	6995	0.448	ng/ul	99
13) Anthracene	17.37	178	5941	0.386	ng/ul	99
15) Fluoranthene	19.29	202	9194	0.462	ng/ul	97
17) Pyrene	19.65	202	10441	0.560	ng/ul#	95
18) Benzo(a)anthracene	21.39	228	9022	0.384	ng/ul	100
19) Chrysene	21.45	228	11102	0.611	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	9558	0.433	ng/ul	98
22) Benzo(k)fluoranthene	23.20	252	10112	0.468	ng/ul#	98
23) Benzo(a)pyrene	23.81	252	9154	0.431	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.57	276	10600	0.436	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.59	278	8675	0.428	ng/ul#	96
26) Benzo(g,h,i)perylene	27.38	276	9211	0.450	ng/ul	97

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

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