

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098050.D
 Acq On : 25 Oct 2018 11:08
 Operator : SJ/JU
 Sample : SST00.161
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.161

Quant Time: Oct 25 13:30:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 25 13:24:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	957	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4732	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	3320	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	8198	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11426	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	9725	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	836	0.11	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	2630	0.10	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	1135	0.100	ng/ul#	91
5) 2-Methylnaphthalene	12.35	142	842	0.111	ng/ul	98
7) Acenaphthylene	14.26	152	1263	0.108	ng/ul	94
8) Acenaphthene	14.60	153	983	0.101	ng/ul	94
9) Fluorene	15.59	166	1160	0.099	ng/ul#	96
12) Phenanthrene	17.33	178	1895	0.097	ng/ul#	94
13) Anthracene	17.43	178	1793	0.108	ng/ul#	92
15) Fluoranthene	19.35	202	2369	0.092	ng/ul	90
17) Pyrene	19.71	202	2407	0.086	ng/ul#	94
18) Benzo(a)anthracene	21.46	228	3660	0.115	ng/ul	97
19) Chrysene	21.52	228	2515	0.082	ng/ul	96
21) Benzo(b)fluoranthene	23.24	252	2618	0.104	ng/ul#	80
22) Benzo(k)fluoranthene	23.29	252	2649	0.093	ng/ul#	78
23) Benzo(a)pyrene	23.91	252	2619	0.114	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	26.72	276	2918	0.112	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.75	278	2456	0.113	ng/ul#	79
26) Benzo(g,h,i)perylene	27.55	276	2437	0.099	ng/ul#	85

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

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