

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025208.D
 Acq On : 03 Mar 2020 11:56
 Operator : CG/JU
 Sample : SST0.480
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.480

Quant Time: Mar 03 12:37:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 03 12:35:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2544	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	9405	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6800	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16985	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	17870	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	17233	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	6403	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	19678	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	10300	0.360	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	7481	0.362	ng/ul	99
7) Acenaphthylene	14.02	152	11322	0.289	ng/ul	95
8) Acenaphthene	14.37	153	8835	0.335	ng/ul	100
9) Fluorene	15.36	166	11232	0.340	ng/ul	96
11) Pentachlorophenol	16.71	266	1510	0.291	ng/ul	99
12) Phenanthrene	17.09	178	19592	0.328	ng/ul	96
13) Anthracene	17.18	178	17174	0.297	ng/ul	96
15) Fluoranthene	19.11	202	24675	0.284	ng/ul	93
17) Pyrene	19.47	202	24281	0.302	ng/ul#	92
18) Benzo(a)anthracene	21.22	228	23623	0.306	ng/ul	98
19) Chrysene	21.27	228	26165	0.340	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	24785	0.367	ng/ul	92
22) Benzo(k)fluoranthene	22.82	252	25796	0.385	ng/ul#	94
23) Benzo(a)pyrene	23.33	252	20660	0.335	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.61	276	26056	0.337	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	21633	0.333	ng/ul#	94
26) Benzo(g,h,i)perylene	26.27	276	22244	0.342	ng/ul#	91

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

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