

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025209.D
 Acq On : 03 Mar 2020 12:32
 Operator : CG/JU
 Sample : SST00.881
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST00.881

Quant Time: Mar 03 13:05:15 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	2186	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	8214	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	5830	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	14222	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	16005	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	15724	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	12714	0.82	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	38565	0.79	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.50	128	20159	0.806	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	14777	0.819	ng/ul	100
7) Acenaphthylene	14.02	152	22339	0.665	ng/ul	94
8) Acenaphthene	14.37	153	17449	0.773	ng/ul	99
9) Fluorene	15.36	166	22061	0.778	ng/ul#	95
11) Pentachlorophenol	16.71	266	3025	0.696	ng/ul	98
12) Phenanthrene	17.09	178	37909	0.758	ng/ul	96
13) Anthracene	17.18	178	34250	0.706	ng/ul	95
15) Fluoranthene	19.11	202	48525	0.668	ng/ul	93
17) Pyrene	19.47	202	47685	0.661	ng/ul#	93
18) Benzo(a)anthracene	21.22	228	48585	0.703	ng/ul	98
19) Chrysene	21.27	228	52633	0.765	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	51646	0.838	ng/ul	90
22) Benzo(k)fluoranthene	22.82	252	53838	0.882	ng/ul#	92
23) Benzo(a)pyrene	23.33	252	43873	0.780	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.61	276	57955	0.822	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	48930	0.826	ng/ul#	91
26) Benzo(g,h,i)perylene	26.27	276	49504	0.833	ng/ul#	89

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

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