

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022420\
 Data File : BM025029.D
 Acq On : 24 Feb 2020 19:35
 Operator : CG/JU
 Sample : SST0.278
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.278

Quant Time: Feb 25 00:51:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 00:46:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1725	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	6803	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	4623	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	11381	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	12793	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	13985	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	2641	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	8018	0.26	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	4204	0.217	ng/ul#	80
5) 2-Methylnaphthalene	12.14	142	3012	0.227	ng/ul	96
7) Acenaphthylene	14.05	152	5696	0.266	ng/ul#	93
8) Acenaphthene	14.39	153	3661	0.204	ng/ul	96
9) Fluorene	15.38	166	4597	0.224	ng/ul#	95
11) Pentachlorophenol	16.72	266	698	0.302	ng/ul	98
12) Phenanthrene	17.11	178	8290	0.229	ng/ul	94
13) Anthracene	17.20	178	8178	0.263	ng/ul	94
15) Fluoranthene	19.13	202	12685	0.324	ng/ul	95
17) Pyrene	19.49	202	12226	0.211	ng/ul#	97
18) Benzo(a)anthracene	21.24	228	11608	0.237	ng/ul	97
19) Chrysene	21.29	228	11549	0.224	ng/ul	97
21) Benzo(b)fluoranthene	22.80	252	11418	0.188	ng/ul	90
22) Benzo(k)fluoranthene	22.84	252	11060	0.178	ng/ul#	89
23) Benzo(a)pyrene	23.36	252	10171	0.191	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.65	276	12846	0.198	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.67	278	10778	0.206	ng/ul#	92
26) Benzo(g,h,i)perylene	26.32	276	10866	0.210	ng/ul#	92

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

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