

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110520\
 Data File : BM027700.D
 Acq On : 05 Nov 2020 01:22
 Operator : JU/CG
 Sample : SST00.103
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST00.1031

Quant Time: Nov 05 04:02:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM110520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 05 03:44:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.40	152	402	0.40	ng/ul	0.00
4) Naphthalene-d8	11.24	136	1422	0.40	ng/ul	0.00
9) Acenaphthene-d10	15.01	164	871	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.72	188	1879	0.40	ng/ul	0.00
19) Chrysene-d12	21.82	240	1572	0.40	ng/ul	0.00
23) Perylene-d12	24.27	264	1427	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/ul	
6) 2-Methylnaphthalene-d10	12.80	152	235	0.10	ng/ul	0.00
17) Fluoranthene-d10	19.71	212	509	0.10	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	11.29	128	398	0.101	ng/ul#	76
7) 2-Methylnaphthalene	12.87	142	278	0.104	ng/ul	98
8) 1-Methylnaphthalene	13.09	142	260	0.103	ng/ul	98
10) Acenaphthylene	14.74	152	392	0.098	ng/ul#	77
11) Acenaphthene	15.07	153	294	0.099	ng/ul	91
12) Fluorene	16.04	166	340	0.098	ng/ul#	97
15) Phenanthrene	17.77	178	558	0.100	ng/ul#	87
16) Anthracene	17.85	178	538	0.100	ng/ul#	87
18) Fluoranthene	19.74	202	648	0.101	ng/ul#	84
20) Pyrene	20.09	202	628	0.098	ng/ul#	84
21) Benzo(a)anthracene	21.80	228	594	0.102	ng/ul#	88
22) Chrysene	21.86	228	605	0.100	ng/ul#	90
24) Benzo(b)fluoranthene	23.52	252	594	0.101	ng/ul#	70
25) Benzo(k)fluoranthene	23.57	252	580	0.099	ng/ul#	76
26) Benzo(a)pyrene	24.16	252	537	0.103	ng/ul#	69
27) Indeno(1,2,3-cd)pyrene	26.80	276	620	0.103	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.82	278	502	0.100	ng/ul#	64
29) Benzo(a,h,i)perylene	27.58	276	521	0.101	ng/ul#	73

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

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