

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011024\
 Data File : BM043809.D
 Acq On : 10 Jan 2024 12:40
 Operator : MA/JU
 Sample : SST0.285
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.2037

Quant Time: Jan 10 14:16:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM011024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 10 14:15:09 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	6660	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	19280	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.606	164	12393	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.358	188	28557	0.400	ng/ul	0.00
17) Chrysene-d12	21.548	240	20439	0.400	ng/ul	0.00
23) Perylene-d12	23.977	264	22607	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.368	96	1619	0.193	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.379	152	5625	0.206	ng/ul	0.01
18) Fluoranthene-d10	19.385	212	13663	0.192	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.402	88	1692	0.199	ng/ul#	78
5) Naphthalene	10.834	128	11233	0.196	ng/ul	96
7) 2-Methylnaphthalene	12.451	142	7406	0.200	ng/ul	98
8) 1-Methylnaphthalene	12.660	142	7568	0.203	ng/ul	98
10) Acenaphthylene	14.328	152	13321	0.196	ng/ul	98
11) Acenaphthene	14.666	153	9795	0.199	ng/ul	100
12) Fluorene	15.665	166	11181	0.199	ng/ul	99
14) Pentachlorophenol	17.021	266	2715	0.408	ng/ul	99
15) Phenanthrene	17.400	178	18151	0.192	ng/ul	98
16) Anthracene	17.498	178	17355	0.201	ng/ul	98
19) Fluoranthene	19.417	202	20864	0.186	ng/ul	100
20) Pyrene	19.780	202	21797	0.194	ng/ul	100
21) Benzo(a)anthracene	21.533	228	17811	0.198	ng/ul	98
22) Chrysene	21.586	228	19635	0.208	ng/ul	99
24) Benzo(b)fluoranthene	23.234	252	18411	0.184	ng/ul	90
25) Benzo(k)fluoranthene	23.284	252	21231	0.203	ng/ul#	87
26) Benzo(a)pyrene	23.874	252	17183	0.202	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.515	276	21641	0.218	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.535	278	16569	0.211	ng/ul#	90
29) Benzo(g,h,i)perylene	27.286	276	18356	0.219	ng/ul	93

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

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