

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022223\
 Data File : BN024089.D
 Acq On : 22 Feb 2023 15:19
 Operator : CG/JU
 Sample : SP6133
 Misc : SFAM SIM SPIKE
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP6133

Quant Time: Feb 23 02:35:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN013023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 17 14:45:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	4595	0.400	ng/ul	0.00
4) Naphthalene-d8	10.634	136	11531	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.479	164	5749	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.230	188	10721	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.426	240	8980	0.400	ng/ul	# 0.00
23) Perylene-d12	23.811	264	6637	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/ul	
6) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng/ul	
18) Fluoranthene-d10	0.000	212	0d	0.000	ng/ul	
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.260	88	18190	3.485	ng/ul#	77
5) Naphthalene	10.678	128	22147	0.702	ng/ul	99
7) 2-Methylnaphthalene	12.300	142	13068	0.637	ng/ul	99
8) 1-Methylnaphthalene	12.520	142	13799	0.657	ng/ul	100
10) Acenaphthylene	14.196	152	15999	0.727	ng/ul	98
11) Acenaphthene	14.543	153	14159	0.732	ng/ul	98
12) Fluorene	15.529	166	15484	0.662	ng/ul	98
14) Pentachlorophenol	16.879	266	4100	1.159	ng/ul	98
15) Phenanthrene	17.272	178	23349	0.730	ng/ul	100
16) Anthracene	17.365	178	18970	0.709	ng/ul	99
19) Fluoranthene	19.290	202	23726	0.795	ng/ul#	96
20) Pyrene	19.653	202	24556	0.794	ng/ul#	93
21) Benzo(a)anthracene	21.408	228	21090	0.691	ng/ul	100
22) Chrysene	21.461	228	24286	0.756	ng/ul	99
24) Benzo(b)fluoranthene	23.086	252	23619	0.806	ng/ul	97
25) Benzo(k)fluoranthene	23.130	252	23659	0.825	ng/ul	98
26) Benzo(a)pyrene	23.703	252	17527	0.718	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.278	276	25853	0.754	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.292	278	20619	0.752	ng/ul	97
29) Benzo(g,h,i)perylene	27.033	276	23030	0.817	ng/ul	96

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

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