

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081523\
 Data File : BN026904.D
 Acq On : 15 Aug 2023 16:55
 Operator : MA/JU
 Sample : 03965-03MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4832MSD

Quant Time: Aug 16 00:54:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN080323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 00:50:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	14705	0.400	ng/ul	0.00
4) Naphthalene-d8	10.938	136	43309	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.735	164	28579	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.481	188	66786	0.400	ng/ul	0.00
17) Chrysene-d12	21.656	240	60844	0.400	ng/ul	0.00
23) Perylene-d12	24.205	264	63810	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.444	96	7418	0.425	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.511	152	21539	0.338	ng/ul	0.00
18) Fluoranthene-d10	19.497	212	61380	0.319	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.477	88	17597	0.987	ng/ul	93
5) Naphthalene	10.988	128	33413	0.273	ng/ul	100
7) 2-Methylnaphthalene	12.583	142	23305	0.314	ng/ul	100
8) 1-Methylnaphthalene	12.797	142	24846	0.320	ng/ul	100
10) Acenaphthylene	14.467	152	37702	0.235	ng/ul	99
11) Acenaphthene	14.800	153	28620	0.258	ng/ul	99
12) Fluorene	15.781	166	33809	0.270	ng/ul	99
15) Phenanthrene	17.523	178	145984	0.665	ng/ul	100
16) Anthracene	17.616	178	56879	0.287	ng/ul	100
19) Fluoranthene	19.529	202	332080	1.314	ng/ul	100
20) Pyrene	19.892	202	303430	1.125	ng/ul	98
21) Benzo(a)anthracene	21.638	228	184607	0.763	ng/ul	100
22) Chrysene	21.694	228	185211	0.775	ng/ul	99
24) Benzo(b)fluoranthene	23.413	252	258560	1.071	ng/ul	98
25) Benzo(k)fluoranthene	23.462	252	131912	0.554	ng/ul	100
26) Benzo(a)pyrene	24.088	252	166704	0.765	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.898	276	163709	0.575	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.928	278	75117	0.331	ng/ul	99
29) Benzo(g,h,i)perylene	27.737	276	127676	0.538	ng/ul	99

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

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