

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023430.D
 Acq On : 25 Dec 2022 03:24
 Operator : CG/JU
 Sample : N6017-01
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBYE8

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/25/2022
 Supervised By :Yogesh Patel 12/25/2022

Quant Time: Dec 25 05:25:06 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Dec 25 05:17:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.990	152	7392	0.400	ng/ul	0.00
4) Naphthalene-d8	10.804	136	24000	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.636	164	14255	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.386	188	32381	0.400	ng/ul	0.00
17) Chrysene-d12	21.572	240	23517	0.400	ng/ul	0.00
23) Perylene-d12	24.019	264	17198	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	30189	3.759	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.394	152	9660	0.260	ng/ul	0.00
18) Fluoranthene-d10	19.416	212	22944	0.270	ng/ul	0.00
Target Compounds						
						Qvalue
10) Acenaphthylene	14.358	152	1288	0.020	ng/ul#	78
15) Phenanthrene	17.428	178	13539	0.128	ng/ul#	86
16) Anthracene	17.521	178	2070	0.024	ng/ul#	81
19) Fluoranthene	19.443	202	37701	0.330	ng/ul	96
20) Pyrene	19.811	202	29417m	0.256	ng/ul	
21) Benzo(a)anthracene	21.555	228	16780	0.183	ng/ul	97
22) Chrysene	21.607	228	19308	0.207	ng/ul	96
24) Benzo(b)fluoranthene	23.267	252	27501m	0.309	ng/ul	
25) Benzo(k)fluoranthene	23.314	252	9268m	0.111	ng/ul	
26) Benzo(a)pyrene	23.908	252	16647	0.236	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.570	276	13407	0.163	ng/ul#	45
28) Dibenzo(a,h)anthracene	26.587	278	3374	0.053	ng/ul#	40
29) Benzo(g,h,i)perylene	27.365	276	13524	0.202	ng/ul#	86

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

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