

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023442.D
 Acq On : 25 Dec 2022 10:33
 Operator : CG/JU
 Sample : N6017-15
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBYG1

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/26/2022
 Supervised By :Sohil Jodhani 12/26/2022

Quant Time: Dec 26 01:22:54 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	7.985	152	6716	0.400	ng/ul	0.00
4) Naphthalene-d8	10.804	136	21999	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.636	164	12938	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.386	188	28796	0.400	ng/ul	0.00
17) Chrysene-d12	21.575	240	20855	0.400	ng/ul	# 0.00
23) Perylene-d12	24.022	264	15307	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	21959	3.009	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.394	152	7831	0.230	ng/ul	0.00
18) Fluoranthene-d10	19.416	212	17782	0.236	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.428	178	4706	0.050	ng/ul	94
19) Fluoranthene	19.443	202	13179	0.130	ng/ul	96
20) Pyrene	19.806	202	10906m	0.107	ng/ul	
21) Benzo(a)anthracene	21.557	228	5749	0.071	ng/ul#	89
22) Chrysene	21.613	228	7253	0.088	ng/ul#	86
24) Benzo(b)fluoranthene	23.276	252	8371m	0.106	ng/ul	
25) Benzo(k)fluoranthene	23.317	252	2631m	0.035	ng/ul	
26) Benzo(a)pyrene	23.914	252	5835	0.093	ng/ul#	60
27) Indeno(1,2,3-cd)pyrene	26.574	276	4039	0.055	ng/ul#	47
28) Dibenzo(a,h)anthracene	26.587	278	1168	0.021	ng/ul#	1
29) Benzo(g,h,i)perylene	27.369	276	5338	0.090	ng/ul#	79

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

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