

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023448.D
 Acq On : 25 Dec 2022 14:10
 Operator : CG/JU
 Sample : N6017-18
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBYG4

Manual Integrations
 APPROVED

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

Quant Time: Dec 26 01:23:46 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	7835	0.400	ng/ul	0.00
4) Naphthalene-d8	10.804	136	25658	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.636	164	15373	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.386	188	34171	0.400	ng/ul	0.00
17) Chrysene-d12	21.575	240	23824	0.400	ng/ul	0.00
23) Perylene-d12	24.022	264	17355	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	30726	3.609	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.394	152	10978	0.277	ng/ul	0.00
18) Fluoranthene-d10	19.416	212	26362	0.306	ng/ul	0.00
Target Compounds						
15) Phenanthrene	17.428	178	10386	0.093	ng/ul	Qvalue 97
19) Fluoranthene	19.443	202	25645	0.222	ng/ul#	67
20) Pyrene	19.811	202	19659m	0.169	ng/ul	
21) Benzo(a)anthracene	21.557	228	10800	0.116	ng/ul#	92
22) Chrysene	21.613	228	14011	0.149	ng/ul#	92
24) Benzo(b)fluoranthene	23.273	252	19405m	0.216	ng/ul	
25) Benzo(k)fluoranthene	23.320	252	6567m	0.078	ng/ul	
26) Benzo(a)pyrene	23.913	252	11332	0.159	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	26.574	276	9614	0.116	ng/ul#	78
28) Dibenzo(a,h)anthracene	26.590	278	2394	0.037	ng/ul#	25
29) Benzo(g,h,i)perylene	27.368	276	10109	0.150	ng/ul#	86

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

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