

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023434.D
 Acq On : 25 Dec 2022 05:47
 Operator : CG/JU
 Sample : N6017-05
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBYF2

Manual Integrations
 APPROVED

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

Quant Time: Dec 26 01:21:44 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	7259	0.400	ng/ul	0.00
4) Naphthalene-d8	10.803	136	23788	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.639	164	14433	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.385	188	31537	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.572	240	19425	0.400	ng/ul	# 0.00
23) Perylene-d12	24.022	264	14500	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	25332	3.212	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.392	152	8640	0.235	ng/ul	0.00
18) Fluoranthene-d10	19.414	212	18565	0.264	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.427	178	6977	0.068	ng/ul#	1
19) Fluoranthene	19.442	202	15946	0.169	ng/ul#	78
20) Pyrene	19.809	202	12492m	0.132	ng/ul	
21) Benzo(a)anthracene	21.554	228	6769	0.090	ng/ul#	80
22) Chrysene	21.610	228	8091	0.105	ng/ul#	72
24) Benzo(b)fluoranthene	23.270	252	11492m	0.153	ng/ul	
25) Benzo(k)fluoranthene	23.314	252	3668m	0.052	ng/ul	
26) Benzo(a)pyrene	23.910	252	6939	0.117	ng/ul#	62
27) Indeno(1,2,3-cd)pyrene	26.574	276	6013	0.087	ng/ul#	45
28) Dibenzo(a,h)anthracene	26.587	278	1509	0.028	ng/ul#	1
29) Benzo(g,h,i)perylene	27.368	276	6057	0.107	ng/ul#	66

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

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