

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023402.D
 Acq On : 24 Dec 2022 10:03
 Operator : CG/JU
 Sample : N6015-06
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXU3

Manual Integrations
 APPROVED

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

Quant Time: Dec 24 23:17:36 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 : Thu Dec 22 02:28:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.990	152	7303	0.400	ng/ul	0.00
4) Naphthalene-d8	10.810	136	24548	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.641	164	14909	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.386	188	34466	0.400	ng/ul	0.00
17) Chrysene-d12	21.581	240	22444	0.400	ng/ul	# 0.00
23) Perylene-d12	24.031	264	16456	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	38269	4.823	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.394	152	12984	0.342	ng/ul	-0.01
18) Fluoranthene-d10	19.416	212	29626	0.365	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.859	128	1474	0.020	ng/ul#	81
15) Phenanthrene	17.428	178	13331	0.119	ng/ul#	85
16) Anthracene	17.521	178	2278	0.024	ng/ul#	83
19) Fluoranthene	19.448	202	35632	0.327	ng/ul#	91
20) Pyrene	19.815	202	26653m	0.243	ng/ul	
21) Benzo(a)anthracene	21.563	228	15811	0.181	ng/ul#	93
22) Chrysene	21.619	228	19917	0.224	ng/ul#	86
24) Benzo(b)fluoranthene	23.279	252	30527m	0.359	ng/ul	
25) Benzo(k)fluoranthene	23.326	252	9639m	0.121	ng/ul	
26) Benzo(a)pyrene	23.922	252	17986	0.267	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.584	276	14751	0.188	ng/ul#	45
28) Dibenzo(a,h)anthracene	26.604	278	3669	0.060	ng/ul#	25
29) Benzo(g,h,i)perylene	27.382	276	14621	0.228	ng/ul#	50

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

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