

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122222\
 Data File : BN023350.D
 Acq On : 22 Dec 2022 21:49
 Operator : CG/JU
 Sample : N6003-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXQ3

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/23/2022
 Supervised By :mohammad ahmed 12/23/2022

Quant Time: Dec 22 23:55:20 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	7.998	152	6487	0.400	ng/ul	0.00
4) Naphthalene-d8	10.815	136	21192	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.645	164	12758	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.390	188	30582	0.400	ng/ul	0.00
17) Chrysene-d12	21.575	240	25359	0.400	ng/ul	0.00
23) Perylene-d12	24.025	264	20750	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.344	96	35990	5.106	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.405	152	10113	0.309	ng/ul	0.00
18) Fluoranthene-d10	19.420	212	27402	0.299	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	14.368	152	1244	0.022	ng/ul#	74
15) Phenanthrene	17.432	178	16158	0.162	ng/ul	99
16) Anthracene	17.525	178	2213	0.027	ng/ul#	85
19) Fluoranthene	19.448	202	50339	0.409	ng/ul	95
20) Pyrene	19.815	202	36272m	0.292	ng/ul	
21) Benzo(a)anthracene	21.560	228	20931	0.212	ng/ul	97
22) Chrysene	21.613	228	25664	0.256	ng/ul	96
24) Benzo(b)fluoranthene	23.276	252	34807m	0.324	ng/ul	
25) Benzo(k)fluoranthene	23.320	252	12236m	0.121	ng/ul	
26) Benzo(a)pyrene	23.919	252	17358	0.204	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.580	276	13346	0.135	ng/ul#	83
28) Dibenzo(a,h)anthracene	26.597	278	3546	0.046	ng/ul#	51
29) Benzo(g,h,i)perylene	27.375	276	9533	0.118	ng/ul#	55

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

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