

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN12222\
 Data File : BN023357.D
 Acq On : 23 Dec 2022 02:08
 Operator : CG/JU
 Sample : N6006-06DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXV9DL

Manual Integrations
 APPROVED

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

Quant Time: Dec 23 03:43:26 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	6289	0.400	ng/ul	0.00
4) Naphthalene-d8	10.815	136	21305	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.645	164	12566	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.390	188	29323	0.400	ng/ul	0.00
17) Chrysene-d12	21.578	240	23255	0.400	ng/ul	0.00
23) Perylene-d12	24.025	264	18248	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.345	96	10241	1.499	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.399	152	3099	0.094	ng/ul	0.00
18) Fluoranthene-d10	19.420	212	7404	0.088	ng/ul	0.00
Target Compounds						
					Qvalue	
11) Acenaphthene	14.706	153	2542	0.054	ng/ul	99
12) Fluorene	15.691	166	4007	0.076	ng/ul	98
15) Phenanthrene	17.432	178	95974	1.005	ng/ul	100
16) Anthracene	17.525	178	21644	0.273	ng/ul	100
19) Fluoranthene	19.448	202	200480	1.777	ng/ul	99
20) Pyrene	19.811	202	134295m	1.181	ng/ul	
21) Benzo(a)anthracene	21.560	228	73111	0.807	ng/ul	99
22) Chrysene	21.616	228	67906	0.738	ng/ul	98
24) Benzo(b)fluoranthene	23.276	252	76995m	0.816	ng/ul	
25) Benzo(k)fluoranthene	23.323	252	25657m	0.290	ng/ul	
26) Benzo(a)pyrene	23.916	252	40644	0.544	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.580	276	26480	0.304	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.597	278	6976	0.103	ng/ul#	71
29) Benzo(g,h,i)perylene	27.375	276	17422	0.245	ng/ul#	93

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

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