

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023378.D
 Acq On : 23 Dec 2022 19:03
 Operator : CG/JU
 Sample : N6008-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBYB1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/24/2022
 Supervised By :Yogesh Patel 12/24/2022

Quant Time: Dec 23 22:44:41 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	7.994	152	6986	0.400	ng/ul	0.00
4) Naphthalene-d8	10.810	136	22619	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.645	164	13432	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.390	188	32806	0.400	ng/ul	0.00
17) Chrysene-d12	21.581	240	25044	0.400	ng/ul	0.00
23) Perylene-d12	24.030	264	19353	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.344	96	10155	1.338	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.399	152	8083	0.231	ng/ul	0.00
18) Fluoranthene-d10	19.420	212	24219	0.267	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.432	178	13108	0.123	ng/ul	96
16) Anthracene	17.525	178	2128	0.024	ng/ul#	85
19) Fluoranthene	19.448	202	45134	0.371	ng/ul	97
20) Pyrene	19.815	202	35108m	0.287	ng/ul	
21) Benzo(a)anthracene	21.563	228	20712	0.212	ng/ul	97
22) Chrysene	21.619	228	24505	0.247	ng/ul	97
24) Benzo(b)fluoranthene	23.279	252	33847m	0.338	ng/ul	
25) Benzo(k)fluoranthene	23.326	252	11143m	0.119	ng/ul	
26) Benzo(a)pyrene	23.922	252	19468	0.246	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.584	276	14755	0.160	ng/ul#	83
28) Dibenzo(a,h)anthracene	26.597	278	3754	0.052	ng/ul#	47
29) Benzo(g,h,i)perylene	27.378	276	14294	0.190	ng/ul#	68

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

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