

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023450.D
 Acq On : 25 Dec 2022 16:05
 Operator : CG/JU
 Sample : N6017-21
 Misc :
 ALS Vial : 79 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBYG9

Quant Time: Dec 26 01:24:04 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	7.985	152	7262	0.400	ng/ul	0.00
4) Naphthalene-d8	10.804	136	23536	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.637	164	13781	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.387	188	30002	0.400	ng/ul	0.00
17) Chrysene-d12	21.576	240	20260	0.400	ng/ul	# 0.00
23) Perylene-d12	24.025	264	14850	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.336	96	20768	2.632	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.393	152	11836	0.325	ng/ul	0.00
18) Fluoranthene-d10	19.416	212	23815	0.325	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	14.359	152	3220	0.053	ng/ul#	91
12) Fluorene	15.687	166	6801	0.118	ng/ul	99
15) Phenanthrene	17.425	178	9841	0.101	ng/ul#	92
16) Anthracene	17.517	178	5570	0.069	ng/ul#	92
19) Fluoranthene	19.444	202	17975	0.183	ng/ul#	89
20) Pyrene	19.807	202	2122	0.021	ng/ul#	48
21) Benzo(a)anthracene	21.561	228	8729	0.111	ng/ul#	80
22) Chrysene	21.614	228	4042	0.050	ng/ul#	47
24) Benzo(b)fluoranthene	23.277	252	3679	0.048	ng/ul#	1
25) Benzo(k)fluoranthene	23.324	252	2202	0.031	ng/ul#	1

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

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