

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110722\
 Data File : BN022559.D
 Acq On : 08 Nov 2022 13:28
 Operator : CG/JU
 Sample : N5407-02
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW4Y2

Quant Time: Nov 09 01:21:51 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 23:39:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.911	152	2225	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.738	136	6616	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.590	164	6167	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.348	188	8300	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.553	240	6797	0.400	ng/ul	# 0.00
23) Perylene-d12	24.000	264	5104	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.224	96	11127	3.848	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.338	152	3833	0.383	ng/ul	0.00
18) Fluoranthene-d10	19.388	212	7306	0.320	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.788	128	7430	0.380	ng/ul#	58
7) 2-Methylnaphthalene	12.410	142	3517	0.286	ng/ul#	77
8) 1-Methylnaphthalene	12.630	142	3099	0.245	ng/ul	90
11) Acenaphthene	14.650	153	1826	0.078	ng/ul	90
12) Fluorene	15.644	166	1321	0.049	ng/ul#	46
14) Pentachlorophenol	17.006	266	1252	0.634	ng/ul	98
15) Phenanthrene	17.390	178	2966	0.106	ng/ul#	10

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

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