

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110722\
 Data File : BN022532.D
 Acq On : 07 Nov 2022 18:23
 Operator : CG/JU
 Sample : N5407-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW4Y2

Quant Time: Nov 07 23:34:53 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 23:30:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.919	152	3289	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.744	136	8704	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.599	164	9983	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.356	188	13086	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.559	240	10797	0.400	ng/ul	# 0.00
23) Perylene-d12	24.014	264	8033	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.224	96	19321	4.520	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	7168	0.544	ng/ul	0.00
18) Fluoranthene-d10	19.392	212	13672	0.377	ng/ul	0.00
Target Compounds				Qvalue		
5) Naphthalene	10.799	128	10401	0.405	ng/ul#	55
7) 2-Methylnaphthalene	12.415	142	5552	0.343	ng/ul#	76
8) 1-Methylnaphthalene	12.635	142	3600	0.216	ng/ul	94
11) Acenaphthene	14.659	153	3832	0.101	ng/ul	89
12) Fluorene	15.654	166	2591	0.059	ng/ul#	45
14) Pentachlorophenol	17.014	266	2332	0.750	ng/ul	99
15) Phenanthrene	17.398	178	5796	0.132	ng/ul#	10

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

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