

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
 Data File : BM036946.D
 Acq On : 11 Oct 2022 15:16
 Operator : CG/JU
 Sample : N4824-04DL 5X
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8P1DL

Quant Time: Oct 12 01:26:50 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 11 09:03:52 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.955	152	8497	0.400 ng/ul	0.00
4) Naphthalene-d8	10.759	136	27952	0.400 ng/ul	# 0.00
9) Acenaphthene-d10	14.579	164	14555	0.400 ng/ul	0.00
13) Phenanthrene-d10	17.317	188	27837	0.400 ng/ul	# 0.00
17) Chrysene-d12	21.485	240	21649	0.400 ng/ul	# 0.00
23) Perylene-d12	23.885	264	21622	0.400 ng/ul	# 0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.352	96	8977	0.749 ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.343	152	2509	0.059 ng/ul	0.00
18) Fluoranthene-d10	19.336	212	4240	0.065 ng/ul	0.00
Target Compounds					
					Qvalue
8) 1-Methylnaphthalene	12.635	142	1035	0.021 ng/ul	99
10) Acenaphthylene	14.302	152	2271	0.038 ng/ul#	18
11) Acenaphthene	14.640	153	49391	0.941 ng/ul	99
12) Fluorene	15.625	166	20182	0.339 ng/ul	99
15) Phenanthrene	17.360	178	6058	0.068 ng/ul#	77
16) Anthracene	17.448	178	14862	0.200 ng/ul#	93
19) Fluoranthene	19.368	202	79231	0.920 ng/ul	98
20) Pyrene	19.726	202	62032	0.691 ng/ul	98
21) Benzo(a)anthracene	21.470	228	10266	0.131 ng/ul#	87
22) Chrysene	21.520	228	8572	0.102 ng/ul#	84
24) Benzo(b)fluoranthene	23.151	252	5169	0.054 ng/ul#	1
26) Benzo(a)pyrene	23.777	252	3050	0.039 ng/ul#	1

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

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