

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112122\
 Data File : BM037672.D
 Acq On : 23 Nov 2022 10:06
 Operator : CG/JU
 Sample : N5591-17
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C9829

Quant Time: Nov 24 01:48:37 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 22:36:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.125	152	8695	0.400	ng/ul	0.00
4) Naphthalene-d8	10.942	136	26519	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.742	164	15929	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.475	188	35235	0.400	ng/ul	0.00
17) Chrysene-d12	21.636	240	24861	0.400	ng/ul	# 0.00
23) Perylene-d12	24.124	264	23650	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.446	96	38020	3.761	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.520	152	8180	0.186	ng/ul	0.00
18) Fluoranthene-d10	19.490	212	19303	0.235	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.997	128	1921	0.024	ng/ul#	87
15) Phenanthrene	17.517	178	5180	0.043	ng/ul#	93
19) Fluoranthene	19.518	202	8446	0.070	ng/ul#	91
20) Pyrene	19.880	202	7051	0.057	ng/ul#	85
21) Benzo(a)anthracene	21.619	228	2986	0.027	ng/ul#	92
22) Chrysene	21.674	228	3629	0.033	ng/ul#	94
24) Benzo(b)fluoranthene	23.361	252	4841	0.045	ng/ul#	5

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

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