

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035528.D
 Acq On : 20 Jun 2022 11:14
 Operator : CG/JU
 Sample : N3226-23
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4P1

Quant Time: Jun 21 02:07:05 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:56:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.829	152	4735	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.616	136	15492	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.455	164	9243	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.204	188	19352	0.400	ng/ul	-0.03
17) Chrysene-d12	21.405	240	13580	0.400	ng/ul	0.00
23) Perylene-d12	23.741	264	14091	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	23575	3.539	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.239	152	5491	0.238	ng/ul	0.00
18) Fluoranthene-d10	19.239	212	14169	0.311	ng/ul	-0.01
Target Compounds						
					Qvalue	
15) Phenanthrene	17.246	178	2586	0.036	ng/ul#	89
21) Benzo(a)anthracene	21.420	228	1167	0.023	ng/ul#	45
22) Chrysene	21.443	228	1327	0.022	ng/ul#	68

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

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