

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035527.D
 Acq On : 20 Jun 2022 10:38
 Operator : CG/JU
 Sample : N3226-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4L5

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/21/2022
 Supervised By :Yogesh Patel 06/25/2022

Quant Time: Jun 21 01:28:14 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	4319	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.622	136	12993	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.459	164	6958	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.208	188	14044	0.400	ng/ul	-0.02
17) Chrysene-d12	21.406	240	11250	0.400	ng/ul	0.00
23) Perylene-d12	23.745	264	11995	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.268	96	11028	1.815	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.233	152	2295	0.119	ng/ul	-0.01
18) Fluoranthene-d10	19.243	212	4009	0.106	ng/ul	-0.01
Target Compounds						
7) 2-Methylnaphthalene	12.310	142	687	0.027	ng/ul	97
8) 1-Methylnaphthalene	12.519	142	577	0.025	ng/ul	99
15) Phenanthrene	17.250	178	2939	0.057	ng/ul#	87
21) Benzo(a)anthracene	21.424	228	1200m	0.028	ng/ul	
22) Chrysene	21.444	228	1572m	0.031	ng/ul	

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

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