

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041222\
 Data File : BM034635.D
 Acq On : 13 Apr 2022 00:09
 Operator : CG/JU
 Sample : N2319-13
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YB2J1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/14/2022

Quant Time: Apr 13 05:10:52 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM041222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 12 15:46:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.884	152	2312	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.687	136	6558	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.528	164	3696	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.270	188	7020m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.481	240	5428	0.400	ng/ul	# 0.01
23) Perylene-d12	23.823	264	2515m	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	18239	4.551	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.309	152	3630	0.418	ng/ul	0.00
18) Fluoranthene-d10	19.298	212	8929	0.565	ng/ul	-0.01
Target Compounds						
2) 1,4-Dioxane	3.290	88	215	0.054	ng/ul#	47
21) Benzo(a)anthracene	21.467	228	299m	0.026	ng/ul	
24) Benzo(b)fluoranthene	23.104	252	544m	0.066	ng/ul	
25) Benzo(k)fluoranthene	23.153	252	332m	0.040	ng/ul	
27) Indeno(1,2,3-cd)pyrene	26.362	276	399	0.032	ng/ul#	64
29) Benzo(g,h,i)perylene	27.100	276	467	0.034	ng/ul#	1

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

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