

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122024\
 Data File : BM049133.D
 Acq On : 21 Dec 2024 11:26
 Operator : RC/JU
 Sample : P5265-01
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E28Y6

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/23/2024
 Supervised By :mohammad ahmed 12/24/2024

Quant Time: Dec 22 23:07:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 18 15:33:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.556	152	8774	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.319	136	27291	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.196	164	15137	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.942	188	32525	0.400	ng/ul	0.00
17) Chrysene-d12	21.143	240	26686	0.400	ng/ul	0.00
23) Perylene-d12	23.300	264	26396m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.151	96	29221	2.888	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.914	152	5860	0.160	ng/ul	-0.01
18) Fluoranthene-d10	18.978	212	10717	0.142	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	16.984	178	4598	0.054	ng/ul#	94
19) Fluoranthene	19.010	202	4035	0.042	ng/ul#	90
20) Pyrene	19.373	202	3466	0.034	ng/ul#	89
22) Chrysene	21.181	228	2033	0.020	ng/ul#	76
24) Benzo(b)fluoranthene	22.666	252	2197m	0.026	ng/ul	
29) Benzo(g,h,i)perylene	26.091	276	1917	0.021	ng/ul#	45

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

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