

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122324\
 Data File : BM049161.D
 Acq On : 24 Dec 2024 00:30
 Operator : RC/JU
 Sample : P5333-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E2910

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/24/2024
 Supervised By :mohammad ahmed 12/26/2024

Quant Time: Dec 24 01:15:51 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	7.552	152	6097	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.309	136	18592	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.187	164	10837	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.938	188	24195m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.140	240	20208m	0.400	ng/ul	0.00
23) Perylene-d12	23.294	264	19201m	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	39498	5.617	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.910	152	8842	0.354	ng/ul	-0.02
18) Fluoranthene-d10	18.974	212	19355	0.338	ng/ul	0.00
Target Compounds						
5) Naphthalene	10.364	128	1075	0.023	ng/ul#	30
7) 2-Methylnaphthalene	11.981	142	1935	0.066	ng/ul	97
8) 1-Methylnaphthalene	12.207	142	1899	0.063	ng/ul	96
15) Phenanthrene	16.976	178	3601m	0.057	ng/ul	
29) Benzo(g,h,i)perylene	26.077	276	2459	0.037	ng/ul#	73

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

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