

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121824\
 Data File : BN035692.D
 Acq On : 19 Dec 2024 04:54
 Operator : RC/JU
 Sample : P5258-12
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E28X6

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/21/2024
 Supervised By :mohammad ahmed 12/24/2024

Quant Time: Dec 19 06:51:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN121624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 16:26:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.276	152	2708	0.400	ng/ul	0.00
4) Naphthalene-d8	10.020	136	7492	0.400	ng/ul	-0.01
9) Acenaphthene-d10	13.939	164	5261	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.709	188	10208	0.400	ng/ul	0.00
17) Chrysene-d12	20.955	240	9134	0.400	ng/ul	0.00
23) Perylene-d12	23.039	264	11898	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.951	96	8236	3.430	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.626	152	2125	0.199	ng/ul	-0.01
18) Fluoranthene-d10	18.763	212	4423	0.166	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	16.751	178	935	0.037	ng/ul#	81
19) Fluoranthene	18.791	202	1400	0.043	ng/ul#	84
20) Pyrene	19.158	202	2212m	0.061	ng/ul	
21) Benzo(a)anthracene	20.941	228	1616	0.053	ng/ul#	83
22) Chrysene	20.990	228	2769	0.088	ng/ul#	85
24) Benzo(b)fluoranthene	22.431	252	1796m	0.051	ng/ul	
26) Benzo(a)pyrene	22.948	252	2019	0.059	ng/ul#	56
27) Indeno(1,2,3-cd)pyrene	25.067	276	1062	0.026	ng/ul#	1
29) Benzo(g,h,i)perylene	25.678	276	1450	0.044	ng/ul#	56

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

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