

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021524\
 Data File : BM044103.D
 Acq On : 15 Feb 2024 16:06
 Operator : MA/JU
 Sample : P1329-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0FLO

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 02/17/2024
 Supervised By :mohammad ahmed 02/19/2024

Quant Time: Feb 15 16:37:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM021524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 15 13:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	5181	0.400	ng/ul	0.00
4) Naphthalene-d8	10.704	136	17907	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.543	164	9847	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.296	188	21698m	0.400	ng/ul	0.00
17) Chrysene-d12	21.497	240	16900m	0.400	ng/ul	0.00
23) Perylene-d12	23.888	264	25532	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	41978	5.721	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.294	152	8329	0.332	ng/ul	0.00
18) Fluoranthene-d10	19.331	212	18196	0.361	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.942	266	337	0.043	ng/ul	92
15) Phenanthrene	17.339	178	3228m	0.047	ng/ul	
19) Fluoranthene	19.359	202	6492	0.084	ng/ul#	93
20) Pyrene	19.721	202	5260	0.065	ng/ul#	86
21) Benzo(a)anthracene	21.482	228	3076	0.040	ng/ul#	92
22) Chrysene	21.535	228	3026	0.036	ng/ul#	91
24) Benzo(b)fluoranthene	23.166	252	4073	0.039	ng/ul#	27
26) Benzo(a)pyrene	23.780	252	2267	0.023	ng/ul#	11

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

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