

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092823\
 Data File : BM042124.D
 Acq On : 28 Sep 2023 20:51
 Operator : MA/JU
 Sample : 04480-13
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBBP0

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/29/2023
 Supervised By :mohammad ahmed 10/03/2023

Quant Time: Sep 29 04:41:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 29 04:39:33 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.942	152	5543	0.400	ng/ul	0.00
4) Naphthalene-d8	10.752	136	14347	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.587	164	11761	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.337	188	17243	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.513	240	10000	0.400	ng/ul	0.00
23) Perylene-d12	23.915	264	9286	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	56837m	7.525	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.330	152	7534	0.419	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	12903	0.313	ng/ul	0.00
Target Compounds						
					Qvalue	
11) Acenaphthene	14.648	153	2378	0.039	ng/ul	84
12) Fluorene	15.633	166	7354	0.108	ng/ul#	69
14) Pentachlorophenol	16.966	266	554	0.061	ng/ul	97
19) Fluoranthene	19.390	202	2928	0.026	ng/ul#	28
20) Pyrene	19.757	202	6654	0.060	ng/ul#	70
22) Chrysene	21.548	228	1945m	0.025	ng/ul	
24) Benzo(b)fluoranthene	23.176	252	1665	0.023	ng/ul#	1

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

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