

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091323\
 Data File : BM041850.D
 Acq On : 14 Sep 2023 03:23
 Operator : MA/JU
 Sample : 04175-06DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EZYX3DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/14/2023
 Supervised By :mohammad ahmed 09/15/2023

Quant Time: Sep 14 05:05:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM091323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 13 16:15:23 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	8304	0.400	ng/ul	0.00
4) Naphthalene-d8	10.785	136	22444	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.615	164	12002	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.367	188	28369	0.400	ng/ul	0.00
17) Chrysene-d12	21.533	240	19086	0.400	ng/ul	0.00
23) Perylene-d12	23.950	264	19663	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	11146	1.055	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.368	152	2605	0.085	ng/ul	0.00
18) Fluoranthene-d10	19.385	212	6130	0.094	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.834	128	31306	0.414	ng/ul	99
15) Phenanthrene	17.409	178	7381	0.060	ng/ul#	95
20) Pyrene	19.780	202	3249	0.021	ng/ul#	83
22) Chrysene	21.571	228	3751m	0.029	ng/ul	
24) Benzo(b)fluoranthene	23.146	252	8685m	0.062	ng/ul	
25) Benzo(k)fluoranthene	23.181	252	7332m	0.057	ng/ul	

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

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