

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080723\
 Data File : BF134654.D
 Acq On : 07 Aug 2023 15:10
 Operator : CG\JU
 Sample : 03887-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MCLEAN-TP4

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/08/2023
 Supervised By :mohammad ahmed 08/08/2023

Quant Time: Aug 08 01:21:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 04 09:29:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.798	152	214069	20.000	ng	0.00
21) Naphthalene-d8	8.075	136	811108	20.000	ng	# 0.00
39) Acenaphthene-d10	9.833	164	382516	20.000	ng	0.00
64) Phenanthrene-d10	11.321	188	606050	20.000	ng	# 0.00
76) Chrysene-d12	13.968	240	446369	20.000	ng	# 0.00
86) Perylene-d12	15.439	264	457710	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1019088	72.365	ng	0.00
7) Phenol-d6	6.434	99	1263976	70.214	ng	0.00
23) Nitrobenzene-d5	7.357	82	731993	56.397	ng	-0.01
42) 2,4,6-Tribromophenol	10.621	330	270524	69.376	ng	0.00
45) 2-Fluorobiphenyl	9.151	172	1246361	54.176	ng	-0.01
79) Terphenyl-d14	12.915	244	1113162	43.963	ng	0.00
Target Compounds						
75) Fluoranthene	12.533	202	111143	3.269	ng	95
78) Pyrene	12.762	202	112616	2.916	ng	99
88) Benzo(b)fluoranthene	15.004	252	62191m	2.228	ng	

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

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