

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090122\
 Data File : BF130093.D
 Acq On : 01 Sep 2022 23:32
 Operator : CG\JU
 Sample : N4423-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-2

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 09/02/2022
 Supervised By :Jagrut Upadhyay 09/02/2022

Quant Time: Sep 02 04:06:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 31 18:15:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.704	152	231607	20.000	ng	0.00
21) Naphthalene-d8	7.981	136	916665	20.000	ng	#-0.01
39) Acenaphthene-d10	9.733	164	478430	20.000	ng	-0.01
64) Phenanthrene-d10	11.216	188	776013	20.000	ng	0.00
76) Chrysene-d12	13.839	240	500929	20.000	ng	#-0.02
86) Perylene-d12	15.245	264	417946	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.316	112	1557479	114.161	ng	0.00
7) Phenol-d6	6.334	99	1777388	104.499	ng	-0.01
23) Nitrobenzene-d5	7.269	82	1253600	87.888	ng	-0.01
42) 2,4,6-Tribromophenol	10.522	330	580949	133.397	ng	-0.01
45) 2-Fluorobiphenyl	9.063	172	2283741	79.141	ng	0.00
79) Terphenyl-d14	12.804	244	2320560	80.333	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	7.681	122	10624m	7.499	ng	

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

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