

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090122\
 Data File : BF130096.D
 Acq On : 02 Sep 2022 01:07
 Operator : CG\JU
 Sample : N4438-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KTS2-WC-S-MH-02-0-20

Quant Time: Sep 02 04:08:11 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	232563	20.000	ng	0.00
21) Naphthalene-d8	7.981	136	909295	20.000	ng	#-0.01
39) Acenaphthene-d10	9.733	164	484686	20.000	ng	-0.01
64) Phenanthrene-d10	11.210	188	778410	20.000	ng	#-0.01
76) Chrysene-d12	13.839	240	495946	20.000	ng	#-0.02
86) Perylene-d12	15.245	264	419559	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.322	112	1556497	113.620	ng	0.01
7) Phenol-d6	6.334	99	1758350	102.955	ng	-0.01
23) Nitrobenzene-d5	7.269	82	1277266	90.273	ng	-0.01
42) 2,4,6-Tribromophenol	10.522	330	575043	130.336	ng	-0.01
45) 2-Fluorobiphenyl	9.063	172	2276694	77.878	ng	0.00
79) Terphenyl-d14	12.804	244	2294989	80.246	ng	0.00
Target Compounds						
32) Benzoic acid	7.698	122	38066	9.921	ng	Qvalue 96

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

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