

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062422\
 Data File : BF128997.D
 Acq On : 24 Jun 2022 18:58
 Operator : CG\JU
 Sample : N3438-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FRAC-TANK-N47935

Quant Time: Jun 25 06:15:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:50:48 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.722	152	98023	20.000	ng	0.00
21) Naphthalene-d8	8.004	136	395132	20.000	ng	0.00
39) Acenaphthene-d10	9.763	164	209179	20.000	ng	0.00
64) Phenanthrene-d10	11.245	188	234178	20.000	ng	0.00
76) Chrysene-d12	13.892	240	160204	20.000	ng	# 0.00
86) Perylene-d12	15.333	264	145246	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.351	112	308040	50.055	ng	0.00
7) Phenol-d6	6.381	99	229619	30.567	ng	0.00
23) Nitrobenzene-d5	7.293	82	672243	79.211	ng	-0.01
42) 2,4,6-Tribromophenol	10.557	330	220488	89.210	ng	0.00
45) 2-Fluorobiphenyl	9.087	172	1232584	82.542	ng	0.00
79) Terphenyl-d14	12.839	244	694288	75.305	ng	0.00

Target Compounds	Qvalue

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

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