

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090222\
 Data File : BF130105.D
 Acq On : 02 Sep 2022 11:56
 Operator : CG\JU
 Sample : N4295-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 XRDS-RECYCLED-SAND-SP-A-COMP

Quant Time: Sep 05 01:16:39 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	213101	20.000	ng	0.00
21) Naphthalene-d8	7.981	136	827866	20.000	ng	-0.01
39) Acenaphthene-d10	9.734	164	427836	20.000	ng	-0.01
64) Phenanthrene-d10	11.216	188	691517	20.000	ng	0.00
76) Chrysene-d12	13.839	240	442512	20.000	ng	#-0.02
86) Perylene-d12	15.245	264	371239	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	1301178	103.657	ng	0.00
7) Phenol-d6	6.328	99	1272114	81.287	ng	-0.02
23) Nitrobenzene-d5	7.269	82	1557621	120.916	ng	-0.01
42) 2,4,6-Tribromophenol	10.522	330	677818	174.045	ng	-0.01
45) 2-Fluorobiphenyl	9.063	172	2649610	102.678	ng	0.00
79) Terphenyl-d14	12.810	244	2662965	104.357	ng	0.00
Target Compounds						
						Qvalue
13) 1,4-Dichlorobenzene	6.722	146	52046	3.383	ng	95
14) 1,2-Dichlorobenzene	6.869	146	38244	2.735	ng	95
30) 1,2,4-Trichlorobenzene	7.922	180	27063	2.193	ng	97

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

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