

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061722\  
 Data File : BF128868.D  
 Acq On : 17 Jun 2022 22:12  
 Operator : CG\JU  
 Sample : N3213-08 10X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SW-W-F1

Quant Time: Jun 18 05:08:24 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF060222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 14 14:53:53 2022  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	77741	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	251900	20.000	ng	# 0.00
39) Acenaphthene-d10	9.827	164	120630	20.000	ng	0.00
64) Phenanthrene-d10	11.321	188	218849	20.000	ng	0.00
76) Chrysene-d12	13.968	240	168977	20.000	ng	# 0.00
86) Perylene-d12	15.427	264	128268	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	12916	2.646	ng	0.01
7) Phenol-d6	6.445	99	13101	2.199	ng	0.00
23) Nitrobenzene-d5	7.345	82	14072	2.601	ng	0.00
42) 2,4,6-Tribromophenol	10.621	330	2314	1.624	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	15887	1.845	ng	0.00
79) Terphenyl-d14	12.904	244	19830	2.039	ng	0.00
Target Compounds						
31) Naphthalene	8.092	128	116561	9.016	ng	# 91
37) 2-Methylnaphthalene	8.786	142	354253	43.082	ng	99
38) 1-Methylnaphthalene	8.886	142	263803	33.206	ng	99
52) Acenaphthene	9.863	154	22463	3.262	ng	# 64
58) Fluorene	10.380	166	56803	7.402	ng	# 91
71) Phenanthrene	11.345	178	184348	16.794	ng	97
78) Pyrene	12.763	202	75045	5.969	ng	# 92
83) Chrysene	13.992	228	26166	2.606	ng	# 45

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

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