

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
 Data File : BF122114.D
 Acq On : 23 Oct 2020 22:55
 Operator : JU/CG
 Sample : L4495-01 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 63202

Manual Integrations
 APPROVED

mohammad
 10/26/2020 10:35:41 AM

Quant Time: Oct 23 23:38:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 12:04:42 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.722	152	103713	20.00	ng	0.00
21) Naphthalene-d8	8.010	136	430466	20.00	ng	0.00
39) Acenaphthene-d10	9.763	164	197371	20.00	ng	0.00
64) Phenanthrene-d10	11.257	188	287114	20.00	ng	0.01
76) Chrysene-d12	13.892	240	265868	20.00	ng	0.00
86) Perylene-d12	15.327	264	261147	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	3524m	0.50	ng	0.08
7) Phenol-d6	6.451	99	7461m	0.82	ng	0.07
23) Nitrobenzene-d5	7.339	82	5366	0.64	ng	0.04
42) 2,4,6-Tribromophenol	10.580	330	712	0.29	ng	0.02
45) 2-Fluorobiphenyl	9.086	172	13012	0.97	ng	0.00
79) Terphenyl-d14	12.833	244	11347	0.81	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	11.280	178	52680	3.35	ng	# 82

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

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