

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
 Data File : BF122111.D
 Acq On : 23 Oct 2020 21:24
 Operator : JU/CG
 Sample : L4487-04 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D-90

Manual Integrations
 APPROVED

mohammad
 10/26/2020 10:35:33 AM

Quant Time: Oct 23 23:37:43 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	94708	20.00	ng	0.00
21) Naphthalene-d8	8.010	136	384097	20.00	ng	0.00
39) Acenaphthene-d10	9.757	164	184711	20.00	ng	0.00
64) Phenanthrene-d10	11.245	188	299259	20.00	ng	0.00
76) Chrysene-d12	13.892	240	226713	20.00	ng	0.00
86) Perylene-d12	15.333	264	232597	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	4078m	0.64	ng	0.09
7) Phenol-d6	6.457	99	7667m	0.93	ng	0.08
23) Nitrobenzene-d5	7.328	82	4795	0.64	ng	0.02
42) 2,4,6-Tribromophenol	10.575	330	828	0.36	ng	0.02
45) 2-Fluorobiphenyl	9.086	172	11654	0.93	ng	0.00
79) Terphenyl-d14	12.833	244	9910	0.83	ng	0.00

Target Compounds	Qvalue

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

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