

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110320\
 Data File : BF122255.D
 Acq On : 3 Nov 2020 22:38
 Operator : JU/CG
 Sample : L4608-01 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OIL-1

Manual Integrations
 APPROVED

mohammad
 11/4/2020 11:01:10 AM

Quant Time: Nov 04 01:08:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 30 10:16:54 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.710	152	74442	20.00	ng	0.00
21) Naphthalene-d8	7.992	136	322373	20.00	ng	0.00
39) Acenaphthene-d10	9.745	164	145310	20.00	ng	0.00
64) Phenanthrene-d10	11.239	188	219541	20.00	ng	# 0.00
76) Chrysene-d12	13.886	240	197112	20.00	ng	0.00
86) Perylene-d12	15.327	264	207887	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	1298m	0.26	ng	0.19
7) Phenol-d6	6.563	99	6010m	0.93	ng	0.19
23) Nitrobenzene-d5	7.363	82	4838m	0.77	ng	0.08
42) 2,4,6-Tribromophenol	10.563	330	902	0.50	ng	0.02
45) 2-Fluorobiphenyl	9.075	172	10445	1.06	ng	0.00
79) Terphenyl-d14	12.821	244	11160	1.08	ng	0.00
Target Compounds						
71) Phenanthrene	11.269	178	34742	2.89	ng	# 72

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

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