

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\
 Data File : BF121500.D
 Acq On : 1 Sep 2020 17:03
 Operator : JU/CG
 Sample : L3859-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-S

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:57:17 PM

Quant Time: Sep 01 18:04:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 09:25:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	281772	20.00	ng	-0.01
21) Naphthalene-d8	8.14	136	1058823	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	542754	20.00	ng	-0.01
64) Phenanthrene-d10	11.38	188	995127	20.00	ng	0.00
76) Chrysene-d12	14.02	240	712586	20.00	ng	0.00
86) Perylene-d12	15.50	264	722961	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1194041	71.10	ng	0.00
7) Phenol-d6	6.47	99	1571350	66.95	ng	-0.02
23) Nitrobenzene-d5	7.42	82	997029	64.38	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	460569	84.01	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1867612	55.52	ng	-0.01
79) Terphenyl-d14	12.97	244	1742366	51.47	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.61	163	202262	5.425	ng	98
71) Phenanthrene	11.40	178	145394	2.847	ng	96
75) Fluoranthene	12.60	202	190948	3.469	ng	95
78) Pyrene	12.82	202	183018	3.596	ng	97
83) Chrysene	14.05	228	89876m	2.065	ng	
88) Benzo(b)fluoranthene	15.06	252	103177m	2.194	ng	

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

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