

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
 Data File : BF118721.D
 Acq On : 27 Jan 2020 16:14
 Operator : CG/JU
 Sample : L1017-07 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-012420

Manual Integrations
 APPROVED

mohammad
 1/28/2020 1:13:11 PM

Quant Time: Jan 27 17:03:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 27 14:24:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	303275	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1067267	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	499148	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	750973	20.00	ng	0.00
76) Chrysene-d12	14.02	240	603354	20.00	ng	0.00
87) Perylene-d12	15.49	264	770576	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	796156	48.76	ng	0.00
7) Phenol-d6	6.48	99	951344	44.82	ng	-0.01
23) Nitrobenzene-d5	7.42	82	590453	30.94	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	222004	38.45	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1148112	37.81	ng	0.00
79) Terphenyl-d14	12.99	244	999180	36.12	ng	0.02
Target Compounds						
50) Dimethylphthalate	9.60	163	105376	3.240	ng	99
71) Phenanthrene	11.40	178	127876	3.530	ng	98
78) Pyrene	12.84	202	91562	2.235	ng	95
81) Benzo(a)anthracene	14.01	228	80406	2.215	ng	# 89
83) Chrysene	14.04	228	78494	2.257	ng	# 95
88) Benzo(b)fluoranthene	15.06	252	98894m	2.042	ng	

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

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