

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030620\
 Data File : BF119270.D
 Acq On : 7 Mar 2020 1:04
 Operator : CG/JU
 Sample : L1762-10 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-08-030520

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:09:35 PM

Quant Time: Mar 07 05:10:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 06 23:38:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	96963	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	374462	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	201951	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	344513	20.00	ng	0.00
76) Chrysene-d12	13.88	240	267201	20.00	ng	0.00
87) Perylene-d12	15.31	264	276103	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	229578	39.57	ng	0.00
7) Phenol-d6	6.37	99	288891	40.94	ng	0.00
23) Nitrobenzene-d5	7.29	82	185988	26.22	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	79348	30.03	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	377357	27.53	ng	0.00
79) Terphenyl-d14	12.82	244	416825	26.86	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.47	163	33441	2.165	ng	98
71) Phenanthrene	11.27	178	78111	4.157	ng	94
75) Fluoranthene	12.46	202	102571	5.143	ng	98
78) Pyrene	12.69	202	89678	4.180	ng	97
81) Benzo(a)anthracene	13.87	228	37723	2.072	ng	96
83) Chrysene	13.91	228	38817	2.140	ng	95
88) Benzo(b)fluoranthene	14.91	252	39917m	2.193	ng	

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

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