

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100720\
 Data File : BF121880.D
 Acq On : 7 Oct 2020 18:31
 Operator : JU/CG
 Sample : L4224-01 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-100620

Manual Integrations
 APPROVED

mohammad
 10/8/2020 12:29:30 PM

Quant Time: Oct 08 00:57:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 14:29:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	130299	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	499327	20.00	ng	-0.01
39) Acenaphthene-d10	9.80	164	223779	20.00	ng	-0.01
64) Phenanthrene-d10	11.29	188	358480	20.00	ng	0.00
76) Chrysene-d12	13.93	240	312034	20.00	ng	0.00
86) Perylene-d12	15.37	264	237503	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	68702	7.84	ng	0.00
7) Phenol-d6	6.41	99	89857	7.81	ng	0.00
23) Nitrobenzene-d5	7.33	82	58345	6.27	ng	-0.01
42) 2,4,6-Tribromophenol	10.60	330	16058	5.77	ng	0.00
45) 2-Fluorobiphenyl	9.12	172	121494	8.22	ng	-0.01
79) Terphenyl-d14	12.87	244	101114	6.56	ng	-0.01
Target Compounds						
71) Phenanthrene	11.31	178	70564	3.613	ng	98
75) Fluoranthene	12.50	202	99259	4.761	ng	99
78) Pyrene	12.73	202	89965	3.946	ng	99
81) Benzo(a)anthracene	13.92	228	44389	2.249	ng	94
83) Chrysene	13.95	228	40638m	2.033	ng	
88) Benzo(b)fluoranthene	14.96	252	38406m	2.419	ng	

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

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