

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
 Data File : BF118417.D
 Acq On : 31 Dec 2019 20:48
 Operator : JU
 Sample : K6463-01RE 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-6-(1-3)RE

Manual Integrations
 APPROVED

mohammad
 1/2/2020 11:50:16 AM

Quant Time: Jan 01 03:12:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	72614	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	273400	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	140538	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	224511	20.00	ng	0.00
76) Chrysene-d12	14.05	240	164908	20.00	ng	0.00
87) Perylene-d12	15.55	264	172810	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	199920	47.00	ng	0.00
7) Phenol-d6	6.49	99	265677	50.35	ng	0.00
23) Nitrobenzene-d5	7.42	82	164334	34.38	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	85050	49.00	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	303271	32.94	ng	0.00
79) Terphenyl-d14	12.99	244	279838	28.15	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.62	163	28399	2.692	ng	Qvalue # 95
71) Phenanthrene	11.42	178	66162	5.406	ng	98
75) Fluoranthene	12.62	202	101624	8.227	ng	99
78) Pyrene	12.85	202	92057	6.836	ng	99
81) Benzo(a)anthracene	14.04	228	44001	3.812	ng	95
83) Chrysene	14.08	228	35513	3.213	ng	96
88) Benzo(b)fluoranthene	15.11	252	48433m	4.222	ng	
90) Benzo(a)pyrene	15.49	252	33685	3.333	ng	96
92) Benzo(g,h,i)perylene	17.54	276	18011	2.126	ng	# 92

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

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