

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
 Data File : BF117622.D
 Acq On : 30 Oct 2019 14:12
 Operator : JU
 Sample : K5542-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 54-54-(0-1.2)

Manual Integrations
 APPROVED

mohammad
 10/31/2019 3:10:36 PM

Quant Time: Oct 31 02:11:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	37523	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	156324	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	73342	20.00	ng	-0.01
64) Phenanthrene-d10	11.25	188	105413	20.00	ng	0.00
76) Chrysene-d12	13.89	240	79556	20.00	ng	0.00
87) Perylene-d12	15.33	264	72376	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	234604	93.02	ng	0.00
7) Phenol-d6	6.39	99	324190	95.70	ng	0.00
23) Nitrobenzene-d5	7.29	82	194080	61.29	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	52024	81.98	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	317650	61.02	ng	-0.01
79) Terphenyl-d14	12.83	244	248729	51.00	ng	0.00
Target Compounds						
10) Phenol	6.40	94	7858	2.273	ng	Qvalue 80
50) Dimethylphthalate	9.48	163	40690	7.782	ng	99
75) Fluoranthene	12.47	202	19992	3.263	ng	97
78) Pyrene	12.70	202	19131	2.800	ng	95
88) Benzo(b)fluoranthene	14.93	252	10741m	2.513	ng	
90) Benzo(a)pyrene	15.28	252	7832	2.016	ng	# 78

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

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