

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102419\
 Data File : BF117515.D
 Acq On : 24 Oct 2019 17:33
 Operator : JU
 Sample : K5499-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 197-16-(0-1)

Manual Integrations
 APPROVED

mohammad
 10/26/2019 10:02:51 AM

Quant Time: Oct 25 03:46:27 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	47652	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	192418	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	97578	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	151830	20.00	ng	0.00
76) Chrysene-d12	13.89	240	91100	20.00	ng	0.00
87) Perylene-d12	15.33	264	100936	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	268106	83.70	ng	0.00
7) Phenol-d6	6.38	99	371725	86.41	ng	0.00
23) Nitrobenzene-d5	7.30	82	222299	57.03	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	56328	66.72	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	282375	40.77	ng	0.00
79) Terphenyl-d14	12.83	244	184019	32.95	ng	0.00
Target Compounds						
10) Phenol	6.39	94	9189	2.093	ng	85
50) Dimethylphthalate	9.48	163	54827	7.882	ng	99
75) Fluoranthene	12.46	202	23777	2.694	ng	99
78) Pyrene	12.69	202	21574	2.757	ng	97
88) Benzo(b)fluoranthene	14.93	252	13725m	2.303	ng	

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

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