

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090519\
 Data File : BF116834.D
 Acq On : 5 Sep 2019 13:35
 Operator : HP/JU
 Sample : K4575-15
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S703A-N-2.0-2.5-082719

Manual Integrations
 APPROVED

mohammad
 9/10/2019 9:52:11 AM

Quant Time: Sep 05 15:08:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	94513	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	383653	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	196014	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	314664	20.00	ng	0.00
76) Chrysene-d12	14.00	240	230259	20.00	ng	0.00
87) Perylene-d12	15.46	264	208631	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	466002	79.88	ng	0.00
7) Phenol-d6	6.48	99	642712	83.90	ng	0.00
23) Nitrobenzene-d5	7.40	82	385357	57.34	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	125605	95.85	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	662570	57.02	ng	0.00
79) Terphenyl-d14	12.94	244	548176	44.04	ng	0.00
Target Compounds						
10) Phenol	6.49	94	17805	2.099	ng	98
50) Dimethylphthalate	9.59	163	67370	4.978	ng	100
71) Phenanthrene	11.39	178	37756	2.155	ng	98
75) Fluoranthene	12.57	202	67947	3.947	ng	99
78) Pyrene	12.80	202	59024	2.884	ng	99
88) Benzo(b)fluoranthene	15.03	252	31410m	2.490	ng	

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

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