

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
 Data File : BF118404.D
 Acq On : 31 Dec 2019 14:40
 Operator : JU
 Sample : K6465-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-10-(10-12)

Manual Integrations
 APPROVED

mohammad
 1/2/2020 11:49:45 AM

Quant Time: Jan 01 00:52:40 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	81000	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	334268	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	190451	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	385355	20.00	ng	0.00
76) Chrysene-d12	14.05	240	253995	20.00	ng	0.00
87) Perylene-d12	15.55	264	229771	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	416893	87.86	ng	0.01
7) Phenol-d6	6.49	99	547497	93.01	ng	0.00
23) Nitrobenzene-d5	7.42	82	331909	56.79	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	192806	81.98	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	602937	48.33	ng	0.00
79) Terphenyl-d14	12.99	244	756146	49.39	ng	0.00
Target Compounds						
10) Phenol	6.50	94	20071	3.034	ng	Qvalue # 74
50) Dimethylphthalate	9.62	163	68040	4.759	ng	99
71) Phenanthrene	11.42	178	101354	4.825	ng	99
75) Fluoranthene	12.62	202	108444	5.115	ng	98
78) Pyrene	12.84	202	126921	6.119	ng	98
81) Benzo(a)anthracene	14.04	228	52975	2.980	ng	98
83) Chrysene	14.08	228	52758m	3.099	ng	
88) Benzo(b)fluoranthene	15.11	252	32057m	2.102	ng	

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

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