

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
 Data File : BF118403.D
 Acq On : 31 Dec 2019 14:12
 Operator : JU
 Sample : K6465-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-11-(1-3)

Manual Integrations
 APPROVED

mohammad
 1/2/2020 11:49:41 AM

Quant Time: Jan 01 00:48:33 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	97307	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	373904	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	210764	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	408429	20.00	ng	0.00
76) Chrysene-d12	14.05	240	283142	20.00	ng	0.00
87) Perylene-d12	15.54	264	257126	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	596313	104.61	ng	0.01
7) Phenol-d6	6.49	99	744354	105.26	ng	0.00
23) Nitrobenzene-d5	7.42	82	455409	69.66	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	262959	101.03	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	866815	62.78	ng	0.00
79) Terphenyl-d14	12.99	244	984931	57.71	ng	0.00
Target Compounds						
10) Phenol	6.50	94	26373	3.318	ng	83
50) Dimethylphthalate	9.62	163	121264	7.665	ng	98
71) Phenanthrene	11.42	178	114084	5.124	ng	98
75) Fluoranthene	12.62	202	161525	7.188	ng	97
78) Pyrene	12.84	202	141944	6.139	ng	99
81) Benzo(a)anthracene	14.04	228	68273	3.445	ng	97
83) Chrysene	14.08	228	60429	3.184	ng	96
88) Benzo(b)fluoranthene	15.11	252	63167m	3.701	ng	
90) Benzo(a)pyrene	15.49	252	41623	2.768	ng	95

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

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