

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
 Data File : BF117588.D
 Acq On : 29 Oct 2019 14:04
 Operator : JU
 Sample : K5503-15
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 172-22-(0-2)

Manual Integrations
 APPROVED

mohammad
 10/30/2019 4:52:55 PM

Quant Time: Oct 30 03:23:18 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	38757	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	148939	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	62335	20.00	ng	-0.01
64) Phenanthrene-d10	11.25	188	82788	20.00	ng	0.00
76) Chrysene-d12	13.89	240	70413	20.00	ng	0.00
87) Perylene-d12	15.33	264	63492	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	240838	92.45	ng	0.00
7) Phenol-d6	6.39	99	340784	97.39	ng	0.00
23) Nitrobenzene-d5	7.29	82	201524	66.80	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	40979	75.98	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	258315	58.39	ng	-0.01
79) Terphenyl-d14	12.83	244	165942	38.44	ng	0.00
Target Compounds						
10) Phenol	6.40	94	10056	2.816	ng	Qvalue 96
50) Dimethylphthalate	9.48	163	33704	7.584	ng	99
71) Phenanthrene	11.27	178	18046	3.754	ng	97
75) Fluoranthene	12.46	202	33211	6.901	ng	94
78) Pyrene	12.69	202	32198	5.324	ng	99
81) Benzo(a)anthracene	13.89	228	17300	3.642	ng	95
83) Chrysene	13.92	228	15981	3.433	ng	97
88) Benzo(b)fluoranthene	14.92	252	17012m	4.537	ng	
90) Benzo(a)pyrene	15.27	252	12041	3.532	ng	# 91

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

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