

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112019\
 Data File : BF117896.D
 Acq On : 20 Nov 2019 16:39
 Operator : JU
 Sample : K5904-19
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 14-(2-4)

Quant Time: Nov 21 03:39:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 18:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	147128	20.00	ng	-0.01
21) Naphthalene-d8	8.20	136	590401	20.00	ng	-0.01
39) Acenaphthene-d10	9.97	164	319854	20.00	ng	0.00
64) Phenanthrene-d10	11.46	188	568177	20.00	ng	-0.01
76) Chrysene-d12	14.10	240	326147	20.00	ng	-0.01
87) Perylene-d12	15.62	264	343482	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	180198	21.68	ng	0.00
7) Phenol-d6	6.53	99	803782	80.17	ng	-0.01
23) Nitrobenzene-d5	7.47	82	601268	60.54	ng	-0.02
42) 2,4,6-Tribromophenol	10.76	330	24601	7.26	ng	-0.01
45) 2-Fluorobiphenyl	9.28	172	1096484	61.50	ng	-0.01
79) Terphenyl-d14	13.04	244	983253	55.10	ng	-0.01
Target Compounds						
10) Phenol	6.55	94	38358	3.682	ng	Qvalue 94
50) Dimethylphthalate	9.67	163	121257	5.406	ng	98
71) Phenanthrene	11.48	178	152388	5.335	ng	97
78) Pyrene	12.90	202	130332	4.945	ng	97
81) Benzo(a)anthracene	14.10	228	58524	2.715	ng	94
83) Chrysene	14.13	228	57163	2.737	ng	97
88) Benzo(b)fluoranthene	15.16	252	72987m	3.271	ng	
90) Benzo(a)pyrene	15.54	252	41889	2.135	ng	# 92

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

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