

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020119\
 Data File : BF112353.D
 Acq On : 1 Feb 2019 16:02
 Operator : JU/SJ
 Sample : K1299-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 0041-FIELD-DUP

Quant Time: Feb 01 23:26:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	153438	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	596399	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	290915	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	532340	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	304268	20.00	ng	-0.01
87) Perylene-d12	15.47	264	323883	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	897679	124.27	ng	0.00
7) Phenol-d6	6.49	99	1098214	109.41	ng	0.00
23) Nitrobenzene-d5	7.40	82	703089	67.93	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	351654	103.85	ng	-0.02
45) 2-Fluorobiphenyl	9.19	172	1296157	63.02	ng	-0.01
79) Terphenyl-d14	12.94	244	1116572	69.12	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	39823	3.616	ng	79
50) Dimethylphthalate	9.58	163	115328	5.102	ng	100
71) Phenanthrene	11.38	178	61216	2.212	ng	97
75) Fluoranthene	12.57	202	69193	2.331	ng	98
78) Pyrene	12.80	202	71545	3.396	ng	98
90) Benzo(a)pyrene	15.40	252	35001	2.008	ng	# 94

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

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