

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021219\
 Data File : BF112519.D
 Acq On : 12 Feb 2019 20:56
 Operator : JU/SJ
 Sample : K1343-23 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-021119-L

Manual Integrations
 APPROVED

Sohil
 2/13/2019 11:12:24 AM

Quant Time: Feb 13 04:57:19 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	124967	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	435962	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	175265	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	316153	20.00	ng	0.00
76) Chrysene-d12	14.02	240	273163	20.00	ng	0.00
87) Perylene-d12	15.50	264	191939	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	371472	63.14	ng	-0.01
7) Phenol-d6	6.49	99	446691	54.64	ng	0.00
23) Nitrobenzene-d5	7.40	82	259580	34.31	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	97101	47.60	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	468324	37.79	ng	-0.01
79) Terphenyl-d14	12.95	244	486255	33.53	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	42371	3.112	ng	Qvalue # 94
71) Phenanthrene	11.39	178	113048	6.878	ng	98
75) Fluoranthene	12.59	202	171444	9.725	ng	99
78) Pyrene	12.82	202	163059	8.622	ng	97
81) Benzo(a)anthracene	14.00	228	68042	4.237	ng	95
83) Chrysene	14.04	228	64959m	4.238	ng	
88) Benzo(b)fluoranthene	15.06	252	60080m	5.234	ng	
90) Benzo(a)pyrene	15.44	252	41928	4.059	ng	# 80
92) Benzo(g,h,i)perylene	17.46	276	23633	3.009	ng	# 82

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

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