

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020819\
 Data File : BF112463.D
 Acq On : 8 Feb 2019 16:53
 Operator : JU/SJ
 Sample : K1409-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-09(0-2)

Manual Integrations
 APPROVED

Sohil
 2/11/2019 10:00:34 AM

Quant Time: Feb 09 05:39:00 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	135255	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	505217	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	230118	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	389268	20.00	ng	-0.01
76) Chrysene-d12	14.01	240	271797	20.00	ng	0.00
87) Perylene-d12	15.50	264	273397	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	708823	111.31	ng	-0.01
7) Phenol-d6	6.49	99	887619	100.32	ng	0.00
23) Nitrobenzene-d5	7.40	82	553706	63.15	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	219365	81.90	ng	-0.02
45) 2-Fluorobiphenyl	9.19	172	1031159	63.38	ng	-0.02
79) Terphenyl-d14	12.94	244	743082	51.49	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	49870	5.137	ng	83
50) Dimethylphthalate	9.58	163	112091	6.269	ng	99
78) Pyrene	12.80	202	46124	2.451	ng	# 92
86) Indeno(1,2,3-cd)pyrene	17.00	276	25108	2.078	ng	96
88) Benzo(b)fluoranthene	15.06	252	43524m	2.662	ng	
90) Benzo(a)pyrene	15.43	252	30429	2.068	ng	# 32
92) Benzo(g,h,i)perylene	17.45	276	33783	3.020	ng	# 84

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

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