

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
 Data File : BF112755.D
 Acq On : 28 Feb 2019 11:59
 Operator : JU/SJ
 Sample : K1627-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-(0-2)

Manual Integrations
 APPROVED

Sohil
 3/1/2019 2:59:46 PM

Quant Time: Mar 01 05:46:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 27 16:08:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	185716	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	656565	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	261376	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	449851	20.00	ng	0.00
76) Chrysene-d12	13.87	240	483787	20.00	ng	0.00
87) Perylene-d12	15.29	264	529327	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	1157738	96.97	ng	0.01
7) Phenol-d6	6.37	99	1421070	94.75	ng	0.00
23) Nitrobenzene-d5	7.30	82	857319	78.13	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	296371	106.81	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1330686	87.80	ng	0.00
79) Terphenyl-d14	12.83	244	1242266	49.78	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	110240	5.720	ng	100
71) Phenanthrene	11.27	178	64822	2.896	ng	99
75) Fluoranthene	12.45	202	100521	4.192	ng	99
78) Pyrene	12.68	202	90863	2.228	ng	98
88) Benzo(b)fluoranthene	14.88	252	73924m	2.159	ng	

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

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