

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
 Data File : BF112812.D
 Acq On : 1 Mar 2019 20:53
 Operator : JU/SJ
 Sample : K1675-37
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Z-3-14-A

Manual Integrations
 APPROVED

Sohil
 3/4/2019 4:55:37 PM

Quant Time: Mar 02 00:35:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	157604	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	542637	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	218483	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	416121	20.00	ng	0.00
76) Chrysene-d12	13.87	240	445215	20.00	ng	0.00
87) Perylene-d12	15.29	264	337631	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	875210	86.38	ng	0.01
7) Phenol-d6	6.37	99	1085471	85.29	ng	0.00
23) Nitrobenzene-d5	7.30	82	628207	69.27	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	250121	107.84	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1047586	77.65	ng	0.00
79) Terphenyl-d14	12.83	244	1128717	49.15	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.49	163	71149	4.416	ng	99
71) Phenanthrene	11.27	178	77934	3.764	ng	97
75) Fluoranthene	12.46	202	174926	7.886	ng	95
78) Pyrene	12.68	202	163324	4.352	ng	98
81) Benzo(a)anthracene	13.86	228	94468	3.062	ng	99
83) Chrysene	13.90	228	84230	2.787	ng	95
88) Benzo(b)fluoranthene	14.89	252	97460m	4.463	ng	
90) Benzo(a)pyrene	15.22	252	55576	2.877	ng	94

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

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