

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030619\
 Data File : BF112912.D
 Acq On : 7 Mar 2019 5:44
 Operator : JU/SJ
 Sample : K1705-13
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EXT-027-SW024-022719

Manual Integrations
 APPROVED

Sohil
 3/7/2019 11:03:09 AM

Quant Time: Mar 07 06:43:34 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	6.73	152	217264	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	686057	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	302770	20.00	ng	-0.01
64) Phenanthrene-d10	11.24	188	591929	20.00	ng	0.00
76) Chrysene-d12	13.86	240	597401	20.00	ng	-0.01
87) Perylene-d12	15.27	264	439518	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1159183	82.99	ng	0.00
7) Phenol-d6	6.37	99	1427147	81.34	ng	0.00
23) Nitrobenzene-d5	7.29	82	846673	73.85	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	346305	107.74	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	1321062	68.74	ng	-0.01
79) Terphenyl-d14	12.82	244	1349897	43.80	ng	-0.01
Target Compounds						
10) Phenol	6.38	94	48308	2.360	ng	88
50) Dimethylphthalate	9.48	163	88694	3.973	ng	98
71) Phenanthrene	11.26	178	222303	7.548	ng	99
72) Anthracene	11.31	178	62418	2.025	ng	96
75) Fluoranthene	12.44	202	325624	10.320	ng	95
78) Pyrene	12.67	202	293060	5.819	ng	98
81) Benzo(a)anthracene	13.85	228	165252	3.991	ng	98
83) Chrysene	13.89	228	149761	3.693	ng	96
88) Benzo(b)fluoranthene	14.87	252	143996m	5.065	ng	
90) Benzo(a)pyrene	15.21	252	88340	3.513	ng	98
92) Benzo(g,h,i)perylene	17.03	276	60529	2.809	ng	97

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

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