

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032919\
 Data File : BF113433.D
 Acq On : 30 Mar 2019 00:40
 Operator : JU/SJ
 Sample : K2153-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CONC-2

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:46:41 PM

Quant Time: Mar 30 03:35:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	139320	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	545209	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	276883	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	575537	20.00	ng	0.00
76) Chrysene-d12	13.84	240	516882	20.00	ng	0.00
87) Perylene-d12	15.25	264	420448	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	157496	18.48	ng	0.00
7) Phenol-d6	6.35	99	766068	71.07	ng	0.00
23) Nitrobenzene-d5	7.27	82	635565	69.19	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	5389	1.58	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1270120	68.47	ng	0.00
79) Terphenyl-d14	12.80	244	1515061	64.61	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.46	163	77446	3.774	ng	99
71) Phenanthrene	11.24	178	222075	7.770	ng	98
72) Anthracene	11.29	178	61602	2.122	ng	98
75) Fluoranthene	12.42	202	285396	8.833	ng	99
78) Pyrene	12.65	202	273138	7.626	ng	98
81) Benzo(a)anthracene	13.83	228	118080	3.991	ng	99
83) Chrysene	13.87	228	108710	3.617	ng	98
86) Indeno(1,2,3-cd)pyrene	16.60	276	52316	2.028	ng	96
88) Benzo(b)fluoranthene	14.85	252	127116m	4.867	ng	
90) Benzo(a)pyrene	15.19	252	89575	3.869	ng	# 97
92) Benzo(g,h,i)perylene	17.02	276	51203	2.536	ng	# 92

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

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