

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030419\
 Data File : BF112830.D
 Acq On : 4 Mar 2019 17:56
 Operator : JU/SJ
 Sample : K1712-11 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Z-2-2

Manual Integrations
 APPROVED

Sohil
 3/6/2019 9:28:19 AM

Quant Time: Mar 05 01:39:20 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	154143	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	554103	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	245807	20.00	ng	0.00
64) Phenanthrene-d10	11.26	188	513447	20.00	ng	0.01
76) Chrysene-d12	13.89	240	453163	20.00	ng	0.01
87) Perylene-d12	15.30	264	333815	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	553915	55.90	ng	0.00
7) Phenol-d6	6.37	99	692858	55.66	ng	0.00
23) Nitrobenzene-d5	7.30	82	392062	42.34	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	192475	73.76	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	639727	36.98	ng	0.00
79) Terphenyl-d14	12.83	244	693344	29.66	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	44199	2.439	ng	98
71) Phenanthrene	11.28	178	526507	20.610	ng	98
73) Carbazole	11.48	167	57671	2.211	ng	97
75) Fluoranthene	12.46	202	780976	28.534	ng	98
78) Pyrene	12.69	202	663967	17.380	ng	99
81) Benzo(a)anthracene	13.87	228	236269	7.523	ng	96
83) Chrysene	13.91	228	330868	10.755	ng	97
86) Indeno(1,2,3-cd)pyrene	16.66	276	139234	5.309	ng	# 92
88) Benzo(b)fluoranthene	14.90	252	360151m	16.680	ng	
89) Benzo(k)fluoranthene	14.92	252	78999m	4.039	ng	
90) Benzo(a)pyrene	15.24	252	176756	9.255	ng	98
92) Benzo(a,h,i)perylene	17.07	276	143701	8.781	ng	# 93

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

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