

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
 Data File : BF112597.D
 Acq On : 18 Feb 2019 13:31
 Operator : JU/SJ
 Sample : K1490-01 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-214

Manual Integrations
 APPROVED

Sohil
 2/19/2019 9:38:34 AM

Quant Time: Feb 18 14:33:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	91993	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	360642	20.00	ng	-0.02
39) Acenaphthene-d10	9.86	164	167244	20.00	ng	-0.02
64) Phenanthrene-d10	11.35	188	292765	20.00	ng	-0.02
76) Chrysene-d12	14.00	240	231742	20.00	ng	-0.01
87) Perylene-d12	15.47	264	249560	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	325074	75.06	ng	-0.02
7) Phenol-d6	6.47	99	415625	69.06	ng	-0.02
23) Nitrobenzene-d5	7.39	82	256127	40.92	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	104328	53.59	ng	-0.02
45) 2-Fluorobiphenyl	9.17	172	388404	32.85	ng	-0.03
79) Terphenyl-d14	12.93	244	320832	26.08	ng	-0.02
Target Compounds						
71) Phenanthrene	11.37	178	88634	5.823	ng	99
75) Fluoranthene	12.57	202	145540	8.915	ng	95
78) Pyrene	12.80	202	132973	8.288	ng	97
81) Benzo(a)anthracene	13.99	228	92314	6.776	ng	97
83) Chrysene	14.02	228	87855	6.756	ng	98
86) Indeno(1,2,3-cd)pyrene	16.97	276	67098	6.512	ng	# 90
88) Benzo(b)fluoranthene	15.04	252	131900m	8.838	ng	
89) Benzo(k)fluoranthene	15.07	252	40979m	2.866	ng	
90) Benzo(a)pyrene	15.42	252	92805	6.911	ng	96
92) Benzo(g,h,i)perylene	17.43	276	64867	6.353	ng	97

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

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