

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030719\
 Data File : BF112934.D
 Acq On : 7 Mar 2019 21:07
 Operator : JU/SJ
 Sample : K1757-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-6AB-S

Manual Integrations
 APPROVED

Sohil
 3/8/2019 9:26:13 AM

Quant Time: Mar 08 01:16:49 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	155658	20.00	ng	-0.01
21) Naphthalene-d8	8.01	136	567983	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	293216	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	586525	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	394514	20.00	ng	-0.02
87) Perylene-d12	15.26	264	416060	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	878722	87.81	ng	-0.01
7) Phenol-d6	6.36	99	1163335	92.55	ng	0.00
23) Nitrobenzene-d5	7.29	82	684440	72.11	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	354830	113.99	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	1265125	67.77	ng	-0.01
79) Terphenyl-d14	12.82	244	1241063	60.98	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.47	163	95270	4.407	ng	99
71) Phenanthrene	11.26	178	97530	3.342	ng	98
75) Fluoranthene	12.44	202	262956	8.411	ng	98
78) Pyrene	12.67	202	214593	6.452	ng	97
81) Benzo(a)anthracene	13.85	228	101662m	3.718	ng	
83) Chrysene	13.89	228	124095	4.633	ng	99
86) Indeno(1,2,3-cd)pyrene	16.61	276	73456m	3.217	ng	
88) Benzo(b)fluoranthene	14.87	252	176126m	6.545	ng	
89) Benzo(k)fluoranthene	14.89	252	66847m	2.742	ng	
90) Benzo(a)pyrene	15.20	252	96552	4.056	ng	99
92) Benzo(g,h,i)perylene	17.02	276	55345	2.713	ng	96

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

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