

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
 Data File : BF112809.D
 Acq On : 1 Mar 2019 19:32
 Operator : JU/SJ
 Sample : K1675-11
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Z-3-1-B

Manual Integrations
 APPROVED

Sohil
 3/4/2019 4:55:34 PM

Quant Time: Mar 02 00:08:07 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	172867	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	609011	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	247252	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	441332	20.00	ng	0.00
76) Chrysene-d12	13.87	240	467520	20.00	ng	0.00
87) Perylene-d12	15.29	264	366273	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	978092	88.01	ng	0.01
7) Phenol-d6	6.37	99	1224057	87.68	ng	0.00
23) Nitrobenzene-d5	7.30	82	718735	70.62	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	285794	108.88	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1214890	80.20	ng	0.00
79) Terphenyl-d14	12.83	244	1235800	51.24	ng	0.00
Target Compounds						
10) Phenol	6.39	94	54175	3.326	ng	92
50) Dimethylphthalate	9.49	163	66987	3.674	ng	96
71) Phenanthrene	11.27	178	148118	6.746	ng	98
75) Fluoranthene	12.46	202	371445	15.789	ng	96
78) Pyrene	12.68	202	349364	8.864	ng	99
81) Benzo(a)anthracene	13.86	228	208845	6.445	ng	100
83) Chrysene	13.90	228	202348	6.375	ng	97
86) Indeno(1,2,3-cd)pyrene	16.64	276	75890	2.805	ng	# 93
88) Benzo(b)fluoranthene	14.89	252	243182m	10.264	ng	
89) Benzo(k)fluoranthene	14.91	252	59293m	2.763	ng	
90) Benzo(a)pyrene	15.23	252	133538	6.372	ng	99
92) Benzo(a,h,i)perylene	17.04	276	68561	3.818	ng	96

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

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