

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032019\
 Data File : BF113173.D
 Acq On : 20 Mar 2019 16:10
 Operator : JU/SJ
 Sample : K1971-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-8(14-15)-031519

Manual Integrations
 APPROVED

Sohil
 3/21/2019 6:36:19 PM

Quant Time: Mar 21 14:42:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 21 12:13:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	107083	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	406984	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	221892	20.00	ng	0.00
64) Phenanthrene-d10	11.23	188	470128	20.00	ng	0.00
76) Chrysene-d12	13.87	240	258547	20.00	ng	0.00
87) Perylene-d12	15.30	264	241244	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	600860	87.28	ng	0.00
7) Phenol-d6	6.37	99	802058	92.75	ng	0.00
23) Nitrobenzene-d5	7.28	82	483138	71.03	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	287418	122.01	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	907030	63.32	ng	0.00
79) Terphenyl-d14	12.82	244	897415	67.29	ng	0.00
Target Compounds						
10) Phenol	6.38	94	59475	5.894	ng	# 77
50) Dimethylphthalate	9.47	163	89093	5.445	ng	# 97
71) Phenanthrene	11.26	178	107573	4.599	ng	97
75) Fluoranthene	12.45	202	92066	3.674	ng	96
78) Pyrene	12.67	202	119097	5.464	ng	97
81) Benzo(a)anthracene	13.86	228	41030	2.290	ng	# 85
83) Chrysene	13.90	228	57544	3.278	ng	# 81
88) Benzo(b)fluoranthene	14.90	252	59099m	3.787	ng	
90) Benzo(a)pyrene	15.24	252	38988	2.825	ng	# 1
92) Benzo(g,h,i)perylene	17.08	276	30132	2.548	ng	# 64

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

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