

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
 Data File : BF112285.D
 Acq On : 28 Jan 2019 20:06
 Operator : JU/SJ
 Sample : K1011-11 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-012519-B

Manual Integrations
 APPROVED

mohammad
 1/29/2019 9:35:19 AM

Quant Time: Jan 29 00:18:49 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	6.84	152	234086	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	883562	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	395866	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	660601	20.00	ng	0.00
76) Chrysene-d12	14.00	240	479420	20.00	ng	0.00
87) Perylene-d12	15.47	264	459120	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	285658	25.92	ng	-0.01
7) Phenol-d6	6.48	99	364779	23.82	ng	-0.01
23) Nitrobenzene-d5	7.40	82	220207	14.36	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	82740	17.96	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	424851	15.18	ng	-0.01
79) Terphenyl-d14	12.94	244	319061	12.53	ng	-0.01
Target Compounds						
71) Phenanthrene	11.39	178	140788	4.099	ng	97
75) Fluoranthene	12.57	202	181049	4.915	ng	98
78) Pyrene	12.80	202	173884	5.239	ng	96
81) Benzo(a)anthracene	13.99	228	75055	2.663	ng	97
83) Chrysene	14.03	228	71570	2.660	ng	97
88) Benzo(b)fluoranthene	15.04	252	84609m	3.081	ng	
90) Benzo(a)pyrene	15.40	252	65056	2.633	ng	# 95

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

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