

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081718\
 Data File : BF108273.D
 Acq On : 18 Aug 2018 5:31
 Operator : JU/SJ
 Sample : J4511-08
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-081518-FL-030

Manual Integrations
 APPROVED

Sohil
 8/20/2018 12:48:46 PM

Quant Time: Aug 18 06:56:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	147863	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	538007	20.00	ng	0.00
38) Acenaphthene-d10	10.06	164	230218	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	368098	20.00	ng	0.00
75) Chrysene-d12	14.21	240	347768	20.00	ng	0.00
86) Perylene-d12	15.78	264	219371	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	903158	110.58	ng	0.02
7) Phenol-d6	6.64	99	1192015	109.83	ng	0.01
23) Nitrobenzene-d5	7.58	82	796952	82.66	ng	0.00
41) 2,4,6-Tribromophenol	10.86	330	237199	103.29	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	1384024	96.52	ng	0.00
78) Terphenyl-d14	13.15	244	1143592	83.61	ng	0.00
Target Compounds						
10) Phenol	6.65	94	28468	2.536	ng	94
49) Dimethylphthalate	9.77	163	147543	9.200	ng	# 98
74) Fluoranthene	12.78	202	55157	2.911	ng	97
77) Pyrene	13.01	202	48428	2.200	ng	99
87) Benzo(b)fluoranthene	15.31	252	28918m	2.206	ng	

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

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