

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110518\
 Data File : BF110476.D
 Acq On : 5 Nov 2018 20:18
 Operator : JU/SJ
 Sample : J5771-01 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-110218-A

Manual Integrations
 APPROVED

Sohil
 11/6/2018 9:01:18 AM

Quant Time: Nov 06 04:36:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	101854	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	407473	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	161725	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	238398	20.00	ng	0.00
76) Chrysene-d12	14.13	240	180516	20.00	ng	0.00
87) Perylene-d12	15.66	264	136660	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	133174	21.29	ng	0.00
7) Phenol-d6	6.57	99	174614	21.83	ng	-0.01
23) Nitrobenzene-d5	7.52	82	106548	15.51	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	28335	17.69	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	186308	18.09	ng	0.00
79) Terphenyl-d14	13.07	244	117128	13.49	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.70	163	27764	2.326	ng	99
71) Phenanthrene	11.51	178	29582	2.371	ng	99
75) Fluoranthene	12.70	202	49218	3.888	ng	96
78) Pyrene	12.94	202	48626	3.757	ng	98
81) Benzo(a)anthracene	14.12	228	26881	2.511	ng	97
83) Chrysene	14.16	228	24612m	2.421	ng	
88) Benzo(b)fluoranthene	15.21	252	22848m	2.895	ng	
90) Benzo(a)pyrene	15.59	252	17405	2.397	ng	# 87

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

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