

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
 Data File : BF111466.D
 Acq On : 15 Dec 2018 14:08
 Operator : JU/SJ
 Sample : J6322-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-2A-C-2-120718

Manual Integrations
 APPROVED

Sohil
 12/17/2018 3:11:55 PM

Quant Time: Dec 16 01:11:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	153856	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	608906	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	273532	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	414829	20.00	ng	0.00
76) Chrysene-d12	13.95	240	299532	20.00	ng	0.00
87) Perylene-d12	15.39	264	293709	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1209419	130.81	ng	0.00
7) Phenol-d6	6.44	99	1579407	128.42	ng	0.00
23) Nitrobenzene-d5	7.36	82	970806	94.94	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	229194	93.54	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1609709	90.90	ng	0.00
79) Terphenyl-d14	12.90	244	1051693	82.45	ng	0.00
Target Compounds						
10) Phenol	6.45	94	94684	6.911	ng	95
50) Dimethylphthalate	9.55	163	132859	6.744	ng	# 96
75) Fluoranthene	12.52	202	56701	2.711	ng	95
78) Pyrene	12.75	202	51551	2.441	ng	95
88) Benzo(b)fluoranthene	14.97	252	37606m	2.196	ng	

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

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