

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111441.D
 Acq On : 15 Dec 2018 1:35
 Operator : JU/SJ
 Sample : J6322-21
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-6-C-2-120718

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:37:13 PM

Quant Time: Dec 15 03:10:48 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	6.80	152	207131	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	757625	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	290395	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	474151	20.00	ng	0.00
76) Chrysene-d12	13.95	240	368696	20.00	ng	0.00
87) Perylene-d12	15.39	264	256568	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1155335	92.82	ng	0.01
7) Phenol-d6	6.44	99	1548374	93.52	ng	0.00
23) Nitrobenzene-d5	7.36	82	1065010	83.71	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	276604	106.33	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1570340	83.53	ng	0.00
79) Terphenyl-d14	12.90	244	1259824	80.24	ng	0.00
Target Compounds						
10) Phenol	6.45	94	82434	4.469	ng	94
50) Dimethylphthalate	9.55	163	120558	5.765	ng	# 96
71) Phenanthrene	11.34	178	89484	3.663	ng	94
75) Fluoranthene	12.53	202	139823	5.850	ng	96
78) Pyrene	12.76	202	118779	4.569	ng	95
81) Benzo(a)anthracene	13.94	228	48596	2.424	ng	99
83) Chrysene	13.98	228	48552	2.299	ng	95
88) Benzo(b)fluoranthene	14.98	252	38130m	2.549	ng	

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

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