

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111566.D
 Acq On : 18 Dec 2018 3:48
 Operator : JU/SJ
 Sample : J6403-61
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-8-B

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:50:03 PM

Quant Time: Dec 18 07:23:28 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	171120	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	670989	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	274155	20.00	ng	-0.01
64) Phenanthrene-d10	11.31	188	400677	20.00	ng	0.00
76) Chrysene-d12	13.94	240	346893	20.00	ng	0.00
87) Perylene-d12	15.38	264	277478	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	738443	71.81	ng	0.00
7) Phenol-d6	6.44	99	944481	69.05	ng	0.00
23) Nitrobenzene-d5	7.35	82	718864	63.80	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	188794	76.88	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1150054	64.80	ng	-0.01
79) Terphenyl-d14	12.89	244	792724	53.66	ng	0.00
Target Compounds						
10) Phenol	6.45	94	51375	3.372	ng	94
50) Dimethylphthalate	9.54	163	177464	8.988	ng	96
75) Fluoranthene	12.52	202	77236	3.824	ng	99
78) Pyrene	12.74	202	70933	2.900	ng	97
81) Benzo(a)anthracene	13.93	228	42528	2.255	ng	97
88) Benzo(b)fluoranthene	14.97	252	49182m	3.040	ng	
90) Benzo(a)pyrene	15.32	252	33472	2.294	ng	94

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

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