

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111565.D
 Acq On : 18 Dec 2018 3:21
 Operator : JU/SJ
 Sample : J6403-57
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-8-A

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:50:00 PM

Quant Time: Dec 18 07:21:38 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	114619	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	452103	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	200129	20.00	ng	-0.01
64) Phenanthrene-d10	11.31	188	298558	20.00	ng	0.00
76) Chrysene-d12	13.94	240	229076	20.00	ng	0.00
87) Perylene-d12	15.38	264	181736	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	569493	82.68	ng	0.00
7) Phenol-d6	6.43	99	728320	79.49	ng	0.00
23) Nitrobenzene-d5	7.35	82	442848	58.33	ng	-0.01
42) 2,4,6-Tribromophenol	10.61	330	123522	68.90	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	762598	58.86	ng	-0.01
79) Terphenyl-d14	12.89	244	530425	54.37	ng	0.00
Target Compounds						
10) Phenol	6.45	94	38828	3.804	ng	Qvalue 93
50) Dimethylphthalate	9.54	163	72613	5.038	ng	97
75) Fluoranthene	12.52	202	59614	3.961	ng	95
78) Pyrene	12.74	202	58432	3.618	ng	96
81) Benzo(a)anthracene	13.93	228	32138	2.580	ng	99
83) Chrysene	13.97	228	30812	2.348	ng	94
88) Benzo(b)fluoranthene	14.97	252	31056m	2.931	ng	
90) Benzo(a)pyrene	15.32	252	22706	2.376	ng	# 92

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

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