

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100818\
 Data File : BF109658.D
 Acq On : 8 Oct 2018 20:05
 Operator : JU/SJ
 Sample : J5270-05 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STOCK-PILE-C

Manual Integrations
 APPROVED

Sohil
 10/9/2018 2:57:35 PM

Quant Time: Oct 09 06:43:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 16:02:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	105664	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	425984	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	185470	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	312444	20.00	ng	0.00
75) Chrysene-d12	13.83	240	231144	20.00	ng	0.00
86) Perylene-d12	15.23	264	203835	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	196858	33.07	ng	0.00
7) Phenol-d6	6.35	99	305810	38.46	ng	-0.01
23) Nitrobenzene-d5	7.26	82	203718	26.36	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	63685	31.51	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	358315	26.61	ng	0.00
78) Terphenyl-d14	12.78	244	266464	22.32	ng	0.00
Target Compounds						
31) Naphthalene	8.00	128	127061	6.450	ng	98
37) 2-Methylnaphthalene	8.69	142	119594	9.268	ng	98
54) Dibenzofuran	9.93	168	39649	2.466	ng	# 89
70) Phenanthrene	11.23	178	105219	6.729	ng	99
74) Fluoranthene	12.41	202	63762	3.903	ng	98
77) Pyrene	12.63	202	57718	3.408	ng	97
80) Benzo(a)anthracene	13.82	228	29949	2.348	ng	# 86
82) Chrysene	13.85	228	38909	2.904	ng	# 89
87) Benzo(b)fluoranthene	14.83	252	28316m	2.514	ng	

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

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