

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100818\
 Data File : BF109661.D
 Acq On : 8 Oct 2018 21:29
 Operator : JU/SJ
 Sample : J5287-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-3

Manual Integrations
 APPROVED

Sohil
 10/9/2018 2:57:41 PM

Quant Time: Oct 09 09:33:54 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	101813	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	386852	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	151564	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	246783	20.00	ng	0.00
75) Chrysene-d12	13.83	240	224492	20.00	ng	0.00
86) Perylene-d12	15.22	264	197797	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	615905	107.38	ng	0.00
7) Phenol-d6	6.35	99	812986	106.10	ng	0.00
23) Nitrobenzene-d5	7.26	82	517537	73.75	ng	0.00
41) 2,4,6-Tribromophenol	10.51	330	150624	91.21	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	772994	70.24	ng	0.00
78) Terphenyl-d14	12.78	244	580577	50.06	ng	0.00
Target Compounds						
10) Phenol	6.36	94	41323	4.703	ng	88
49) Dimethylphthalate	9.45	163	82434	7.416	ng	99
74) Fluoranthene	12.41	202	43231	3.351	ng	98
77) Pyrene	12.63	202	43111	2.621	ng	100
80) Benzo(a)anthracene	13.82	228	25134m	2.029	ng	
82) Chrysene	13.85	228	29979m	2.304	ng	
87) Benzo(b)fluoranthene	14.83	252	34507m	3.157	ng	

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

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