

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
 Data File : BF109905.D
 Acq On : 16 Oct 2018 15:13
 Operator : JU/SJ
 Sample : J5526-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-4B-C-101518-2

Manual Integrations
 APPROVED

Sohil
 10/17/2018 11:15:17 AM

Quant Time: Oct 16 16:38:59 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 14:58:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	267527	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	976482	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	365653	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	547104	20.00	ng	0.00
76) Chrysene-d12	14.16	240	634306	20.00	ng	0.00
87) Perylene-d12	15.69	264	665754	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1284944	76.13	ng	0.00
7) Phenol-d6	6.60	99	1584260	75.60	ng	0.00
23) Nitrobenzene-d5	7.54	82	957231	66.04	ng	-0.01
42) 2,4,6-Tribromophenol	10.81	330	234687	74.13	ng	0.00
45) 2-Fluorobiphenyl	9.34	172	1458609	63.65	ng	0.00
79) Terphenyl-d14	13.10	244	1145182	37.91	ng	0.00
Target Compounds						
10) Phenol	6.61	94	74642	3.299	ng	Qvalue # 59
50) Dimethylphthalate	9.73	163	205564	7.939	ng	98
71) Phenanthrene	11.53	178	86911	3.068	ng	100
75) Fluoranthene	12.73	202	201601	7.125	ng	98
78) Pyrene	12.96	202	188717	4.051	ng	99
81) Benzo(a)anthracene	14.14	228	107366	2.874	ng	99
83) Chrysene	14.18	228	102530	2.883	ng	96
88) Benzo(b)fluoranthene	15.23	252	146189m	3.809	ng	
90) Benzo(a)pyrene	15.62	252	106466	3.016	ng	# 97

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

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