

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
 Data File : BF109711.D
 Acq On : 9 Oct 2018 22:47
 Operator : JU/SJ
 Sample : J5326-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FL-PIT-3

Manual Integrations
 APPROVED

Sohil
 10/10/2018 11:37:18 AM

Quant Time: Oct 10 10:18:03 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	81494	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	308111	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	124361	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	201078	20.00	ng	0.00
75) Chrysene-d12	13.83	240	206934	20.00	ng	0.00
86) Perylene-d12	15.23	264	187317	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	671606	146.28	ng	0.00
7) Phenol-d6	6.35	99	924082	150.67	ng	0.00
23) Nitrobenzene-d5	7.26	82	594054	106.28	ng	0.00
41) 2,4,6-Tribromophenol	10.51	330	170022	125.47	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	958775	106.18	ng	0.00
78) Terphenyl-d14	12.78	244	749470	70.11	ng	0.00
Target Compounds						
10) Phenol	6.36	94	33754	4.799	ng	87
49) Dimethylphthalate	9.45	163	93003	10.197	ng	100
70) Phenanthrene	11.23	178	28308	2.813	ng	98
74) Fluoranthene	12.41	202	82083	7.808	ng	99
77) Pyrene	12.63	202	82674	5.453	ng	98
80) Benzo(a)anthracene	13.82	228	72469	6.346	ng	98
82) Chrysene	13.85	228	53033	4.421	ng	96
85) Indeno(1,2,3-cd)pyrene	16.57	276	23967	2.792	ng	94
87) Benzo(b)fluoranthene	14.83	252	84975m	8.209	ng	
88) Benzo(k)fluoranthene	14.86	252	30603m	2.942	ng	
89) Benzo(a)pyrene	15.17	252	50198	5.344	ng	99
91) Benzo(a,h,i)perylene	16.97	276	21679	2.814	ng	98

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

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