

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
 Data File : BF110933.D
 Acq On : 21 Nov 2018 17:28
 Operator : JU/SJ
 Sample : J5772-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-112018-B

Manual Integrations
 APPROVED

mohammad
 11/23/2018 11:49:09 AM

Quant Time: Nov 21 23:11:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	61845	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	247468	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	105471	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	161114	20.00	ng	0.00
76) Chrysene-d12	14.13	240	134992	20.00	ng	0.00
87) Perylene-d12	15.67	264	95280	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	344722	90.54	ng	0.00
7) Phenol-d6	6.57	99	429855	88.29	ng	0.00
23) Nitrobenzene-d5	7.51	82	277204	64.51	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	78226	73.61	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	470841	63.76	ng	0.00
79) Terphenyl-d14	13.06	244	318253	44.15	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	21833	3.867	ng	Qvalue # 76
50) Dimethylphthalate	9.69	163	56067	7.124	ng	99
75) Fluoranthene	12.70	202	33334	3.852	ng	98
78) Pyrene	12.93	202	34116	3.282	ng	98
81) Benzo(a)anthracene	14.12	228	21491m	2.664	ng	
83) Chrysene	14.16	228	21571	2.806	ng	99
88) Benzo(b)fluoranthene	15.22	252	23347m	4.096	ng	
90) Benzo(a)pyrene	15.60	252	14347	2.770	ng	# 92

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

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