

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021521\
 Data File : BF123209.D
 Acq On : 15 Feb 2021 17:28
 Operator : JU/CG
 Sample : M1402-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-N9994

Manual Integrations
 APPROVED

mohammad
 2/16/2021 10:44:49 AM

Quant Time: Feb 15 23:44:38 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF021221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 12 16:59:49 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	146048	20.00	ng	0.00
21) Naphthalene-d8	8.151	136	576116	20.00	ng	0.00
39) Acenaphthene-d10	9.910	164	284735	20.00	ng	0.00
64) Phenanthrene-d10	11.398	188	538176	20.00	ng	0.00
76) Chrysene-d12	14.051	240	398385	20.00	ng	0.00
86) Perylene-d12	15.527	264	312259	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	991757	102.93	ng	0.00
7) Phenol-d6	6.492	99	1222564	96.08	ng	0.00
23) Nitrobenzene-d5	7.428	82	776471	70.82	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	340522	110.74	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1352949	71.01	ng	0.00
79) Terphenyl-d14	12.992	244	1371051	68.97	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.504	94	75356	5.39	ng	90
50) Dimethylphthalate	9.622	163	94890	4.60	ng	100
71) Phenanthrene	11.422	178	67652	2.13	ng	97
75) Fluoranthene	12.616	202	227599	7.14	ng	99
78) Pyrene	12.845	202	190913	6.48	ng	97
81) Benzo(a)anthracene	14.039	228	58398	2.18	ng	98
83) Chrysene	14.074	228	71373m	2.70	ng	
88) Benzo(b)fluoranthene	15.092	252	56558m	2.58	ng	

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

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