

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
 Data File : BF125947.D
 Acq On : 26 Oct 2021 22:15
 Operator : JU/CG
 Sample : M4348-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 358

Manual Integrations
 APPROVED

mohammad
 10/27/2021 11:49:02 AM

Quant Time: Oct 27 09:04:30 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 25 18:49:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	163167	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	638561	20.000	ng	#-0.01
39) Acenaphthene-d10	9.904	164	337225	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	530270	20.000	ng	0.00
76) Chrysene-d12	14.021	240	329069	20.000	ng	-0.01
86) Perylene-d12	15.492	264	374548	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	1009081	92.117	ng	0.00
7) Phenol-d6	6.486	99	1278390	85.086	ng	0.00
23) Nitrobenzene-d5	7.428	82	834705	67.520	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	254798	86.674	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1164424	61.063	ng	0.00
79) Terphenyl-d14	12.974	244	849004	42.202	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.410	178	95497	3.471	ng	98
75) Fluoranthene	12.598	202	122024	4.399	ng	98
78) Pyrene	12.821	202	98414	3.214	ng	96
84) Bis(2-ethylhexyl)phtha...	14.015	149	54938	5.091	ng	94
88) Benzo(b)fluoranthene	15.062	252	53594m	2.168	ng	

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

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