

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061521\
 Data File : BF124333.D
 Acq On : 15 Jun 2021 20:51
 Operator : JU/CG
 Sample : M2674-07 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TMR-DS-01-060921

Manual Integrations
 APPROVED

mohammad
 6/16/2021 3:53:49 PM

Quant Time: Jun 15 21:16:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 15 20:09:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	36415	20.00	ng	0.00
21) Naphthalene-d8	8.069	136	136527	20.00	ng	# 0.00
39) Acenaphthene-d10	9.822	164	79709	20.00	ng	0.00
64) Phenanthrene-d10	11.310	188	150467	20.00	ng	0.00
76) Chrysene-d12	13.957	240	95249	20.00	ng	# 0.00
86) Perylene-d12	15.409	264	77070	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	50958	21.07	ng	0.00
7) Phenol-d6	6.439	99	57845	17.79	ng	0.00
23) Nitrobenzene-d5	7.345	82	56058	16.28	ng	-0.01
42) 2,4,6-Tribromophenol	10.616	330	24691	21.86	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	105797	16.70	ng	0.00
79) Terphenyl-d14	12.898	244	108514	15.64	ng	0.00
Target Compounds						
10) Phenol	6.451	94	22829	6.50	ng	# 74
17) 2-Methylphenol	7.057	107	33511m	15.31	ng	
20) 3+4-Methylphenols	7.216	107	76484	26.38	ng	# 70
27) 2,4-Dimethylphenol	7.734	122	61500	28.27	ng	95

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

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