

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070121\
 Data File : BF124512.D
 Acq On : 1 Jul 2021 15:27
 Operator : JU/CG
 Sample : M2896-03DL 50X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FS-02DL

Manual Integrations
 APPROVED

mohammad
 7/2/2021 12:02:04 PM

Quant Time: Jul 01 15:53:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 01 14:15:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.681	152	31162	20.000	ng	0.00
21) Naphthalene-d8	7.963	136	112542	20.000	ng	# 0.00
39) Acenaphthene-d10	9.716	164	65697	20.000	ng	0.00
64) Phenanthrene-d10	11.204	188	126795	20.000	ng	# 0.00
76) Chrysene-d12	13.851	240	73350	20.000	ng	0.00
86) Perylene-d12	15.280	264	73378	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	0.000	112	0d	0.000	ng	
7) Phenol-d6	6.351	99	1590	0.615	ng	0.01
23) Nitrobenzene-d5	7.257	82	1669	0.615	ng	0.00
42) 2,4,6-Tribromophenol	10.516	330	950	1.210	ng	0.00
45) 2-Fluorobiphenyl	9.039	172	3155	0.588	ng	0.00
79) Terphenyl-d14	12.786	244	7484	1.528	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.357	94	24426	8.732	ng	89
17) 2-Methylphenol	6.957	107	21741m	12.338	ng	
20) 3+4-Methylphenols	7.116	107	45651	19.848	ng	# 72
27) 2,4-Dimethylphenol	7.634	122	30464	17.328	ng	# 89

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

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