

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063021\
 Data File : BF124504.D
 Acq On : 30 Jun 2021 21:38
 Operator : JU/CG
 Sample : M2896-04 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FS-02

Manual Integrations
 APPROVED

mohammad
 7/1/2021 2:52:06 PM

Quant Time: Jul 01 05:16:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 29 12:49:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.692	152	35576	20.000	ng	0.00
21) Naphthalene-d8	7.986	136	171930	20.000	ng	# 0.01
39) Acenaphthene-d10	9.739	164	86421	20.000	ng	# 0.01
64) Phenanthrene-d10	11.222	188	158078	20.000	ng	0.00
76) Chrysene-d12	13.857	240	86514	20.000	ng	0.00
86) Perylene-d12	15.286	264	93363	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	30815m	13.622	ng	0.05
7) Phenol-d6	6.410	99	38534	13.046	ng	0.06
23) Nitrobenzene-d5	7.275	82	140175	33.805	ng	0.01
42) 2,4,6-Tribromophenol	10.539	330	15245	14.765	ng	0.02
45) 2-Fluorobiphenyl	9.069	172	56503	8.009	ng	0.02
79) Terphenyl-d14	12.798	244	61854	10.709	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.428	94	833533	260.996	ng	97
17) 2-Methylphenol	7.028	107	641847	319.054	ng	95
20) 3+4-Methylphenols	7.186	107	1306351	497.489	ng	# 68
22) Acetophenone	7.122	105	53880	11.258	ng	# 85
27) 2,4-Dimethylphenol	7.698	122	565150	210.424	ng	94

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

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