

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070121\
 Data File : BF124514.D
 Acq On : 1 Jul 2021 16:34
 Operator : JU/CG
 Sample : M2896-04DL 100X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FS-02DL

Manual Integrations
 APPROVED

mohammad
 7/2/2021 12:02:06 PM

Quant Time: Jul 01 17:02:12 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	30228	20.000	ng	0.00
21) Naphthalene-d8	7.963	136	117260	20.000	ng	# 0.00
39) Acenaphthene-d10	9.716	164	68080	20.000	ng	0.00
64) Phenanthrene-d10	11.204	188	130539	20.000	ng	# 0.00
76) Chrysene-d12	13.851	240	75082	20.000	ng	0.00
86) Perylene-d12	15.280	264	69963	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	867	0.451	ng	0.06
7) Phenol-d6	6.345	99	2839	1.131	ng	0.00
23) Nitrobenzene-d5	7.228	82	12105m	4.280	ng	-0.03
42) 2,4,6-Tribromophenol	10.516	330	1012	1.244	ng	0.00
45) 2-Fluorobiphenyl	9.039	172	4852	0.873	ng	0.00
79) Terphenyl-d14	12.786	244	8367	1.669	ng	0.00
Target Compounds						
10) Phenol	6.357	94	73811	27.201	ng	Qvalue 95
17) 2-Methylphenol	6.957	107	52035	30.442	ng	# 92
20) 3+4-Methylphenols	7.122	107	111534	49.990	ng	# 70
27) 2,4-Dimethylphenol	7.639	122	46236	25.241	ng	91

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

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