

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031021\
 Data File : BM029080.D
 Acq On : 10 Mar 2021 21:26
 Operator : CG/JU
 Sample : M1603-03RE
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 33-NORTHRE

Manual Integrations
 APPROVED

mohammad
 3/11/2021 9:26:25 AM

Quant Time: Mar 10 23:40:50 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 09 10:08:02 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	11595	20.00	ng	0.00
21) Naphthalene-d8	10.457	136	45352	20.00	ng	# 0.00
39) Acenaphthene-d10	14.333	164	26851	20.00	ng	0.00
64) Phenanthrene-d10	17.080	188	54895	20.00	ng	0.00
76) Chrysene-d12	21.262	240	54528	20.00	ng	# 0.00
86) Perylene-d12	23.439	264	52356	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.240	112	10097	14.43	ng	0.00
7) Phenol-d6	6.857	99	52325	56.63	ng	-0.01
23) Nitrobenzene-d5	8.834	82	40879	48.63	ng	-0.01
42) 2,4,6-Tribromophenol	15.833	330	720	2.26	ng	0.00
45) 2-Fluorobiphenyl	12.945	172	100448	49.80	ng	0.00
79) Terphenyl-d14	19.709	244	144871	49.06	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.887	94	2877	3.13	ng	91
50) Dimethylphthalate	13.798	163	5849	2.87	ng	# 95
71) Phenanthrene	17.121	178	6989	2.14	ng	# 93
75) Fluoranthene	19.139	202	11604	3.23	ng	78
78) Pyrene	19.504	202	12430	3.10	ng	# 81
88) Benzo(b)fluoranthene	22.792	252	7989m	2.16	ng	

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

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