

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM043021\
 Data File : BM029722.D
 Acq On : 29 Apr 2021 23:03
 Operator : CG/JU
 Sample : M2109-01DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PE-1DL

Quant Time: Apr 30 01:09:29 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM043021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 29 17:34:54 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.598	152	70640	20.00	ng	0.00
21) Naphthalene-d8	10.375	136	252061	20.00	ng	0.00
39) Acenaphthene-d10	14.251	164	167972	20.00	ng	0.00
64) Phenanthrene-d10	16.998	188	364253	20.00	ng	0.00
76) Chrysene-d12	21.180	240	404297	20.00	ng	0.00
86) Perylene-d12	23.356	264	433755	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.198	112	62790	18.53	ng	0.00
7) Phenol-d6	6.781	99	72579	15.78	ng	0.00
23) Nitrobenzene-d5	8.751	82	49668	11.39	ng	0.00
42) 2,4,6-Tribromophenol	15.745	330	38777	14.54	ng	0.00
45) 2-Fluorobiphenyl	12.863	172	159714	12.33	ng	0.00
79) Terphenyl-d14	19.627	244	256230	11.63	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.428	128	61246	4.75	ng	# 96
37) 2-Methylnaphthalene	12.051	142	234153	23.62	ng	98
38) 1-Methylnaphthalene	12.269	142	142395	15.08	ng	97
46) 1,1'-Biphenyl	13.074	154	39435	3.11	ng	96
58) Fluorene	15.298	166	28913	2.18	ng	# 83
71) Phenanthrene	17.039	178	109552	5.34	ng	96

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

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