

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121720\
 Data File : BM028149.D
 Acq On : 17 Dec 2020 19:57
 Operator : JU/CG
 Sample : L5113-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CS1

Quant Time: Dec 18 00:21:25 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 16 01:58:38 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.210	152	134228	20.00	ng	0.00
21) Naphthalene-d8	11.039	136	577828	20.00	ng	# 0.00
39) Acenaphthene-d10	14.833	164	353395	20.00	ng	-0.01
64) Phenanthrene-d10	17.557	188	811933	20.00	ng	# 0.00
76) Chrysene-d12	21.668	240	774894	20.00	ng	# 0.00
86) Perylene-d12	24.033	264	661474	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.704	112	500657	61.46	ng	0.00
7) Phenol-d6	7.351	99	1071485	92.12	ng	0.00
23) Nitrobenzene-d5	9.387	82	765510	65.74	ng	-0.01
42) 2,4,6-Tribromophenol	16.310	330	75362	16.06	ng	0.00
45) 2-Fluorobiphenyl	13.463	172	1810023	66.59	ng	-0.01
79) Terphenyl-d14	20.127	244	2867526	67.15	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.375	94	33488	3.03	ng	88
50) Dimethylphthalate	14.280	163	195406	6.66	ng	100
78) Pyrene	19.939	202	117502	2.09	ng	98
88) Benzo(b)fluoranthene	23.315	252	120486	2.67	ng	# 86
90) Benzo(a)pyrene	23.927	252	87681	2.17	ng	# 89
92) Benzo(g,h,i)perylene	27.209	276	71590	2.04	ng	# 85

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

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