

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032421\
 Data File : BP005036.D
 Acq On : 24 Mar 2021 16:24
 Operator : CG/JU
 Sample : M1557-06DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 24 17:05:15 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 24 10:16:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	27299	20.00	ng	0.00
21) Naphthalene-d8	10.516	136	103161	20.00	ng	# 0.00
39) Acenaphthene-d10	14.381	164	66923	20.00	ng	0.00
64) Phenanthrene-d10	17.139	188	148940	20.00	ng	# 0.00
76) Chrysene-d12	21.233	240	149235	20.00	ng	0.00
86) Perylene-d12	23.521	264	152537	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.317	112	53613	30.22	ng	0.00
7) Phenol-d6	6.899	99	68122	26.81	ng	0.00
23) Nitrobenzene-d5	8.887	82	46198	19.21	ng	0.00
42) 2,4,6-Tribromophenol	15.875	330	25809	29.60	ng	0.00
45) 2-Fluorobiphenyl	12.993	172	99848	20.18	ng	0.00
79) Terphenyl-d14	19.710	244	169478	20.67	ng	0.00
Target Compounds						
						Qvalue
8) 2-Chlorophenol	7.293	128	19046	10.18	ng	# 80
10) Phenol	6.928	94	62680	24.07	ng	81
17) 2-Methylphenol	8.169	107	51895	31.23	ng	# 89
26) 2-Nitrophenol	9.634	139	29699	30.22	ng	# 59
27) 2,4-Dimethylphenol	9.699	122	20336	12.87	ng	# 82
29) 2,4-Dichlorophenol	10.163	162	22439	12.69	ng	93
36) 4-Chloro-3-methylphenol	11.810	107	34485	16.14	ng	# 78
43) 2,4,6-Trichlorophenol	12.804	196	11773	7.82	ng	95
44) 2,4,5-Trichlorophenol	12.875	196	26348	16.15	ng	# 93
54) 2,4-Dinitrophenol	14.498	184	13977	19.00	ng	# 74
56) 4-Nitrophenol	14.604	139	24547	26.76	ng	# 34
65) 4,6-Dinitro-2-methylph...	15.516	198	29893	27.50	ng	86
70) Pentachlorophenol	16.787	266	33419	26.08	ng	96

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

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