

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081320\
 Data File : BP002970.D
 Acq On : 13 Aug 2020 17:30
 Operator : CG/JU
 Sample : L3648-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BS-STOCK-4

Quant Time: Aug 13 18:53:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 11 17:14:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	354253	20.00	ng	-0.01
21) Naphthalene-d8	10.47	136	1470222	20.00	ng	-0.01
39) Acenaphthene-d10	14.33	164	887414	20.00	ng	-0.01
64) Phenanthrene-d10	17.08	188	1855766	20.00	ng	-0.01
76) Chrysene-d12	21.17	240	1824547	20.00	ng	0.00
86) Perylene-d12	23.42	264	1930112	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.30	112	1434961	62.00	ng	0.00
7) Phenol-d6	6.86	99	1240994	38.62	ng	-0.02
23) Nitrobenzene-d5	8.83	82	1950478	71.85	ng	-0.01
42) 2,4,6-Tribromophenol	15.82	330	1340395	121.53	ng	-0.01
45) 2-Fluorobiphenyl	12.95	172	5393158	83.04	ng	-0.01
79) Terphenyl-d14	19.66	244	6946300	76.22	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
26) 2-Nitrophenol	9.59	139	30	4.523	ng	# 1
32) Benzoic acid	9.93	122	143	7.695	ng	96
48) 2-Nitroaniline	13.39	65	80	3.977	ng	# 10
51) 2,6-Dinitrotoluene	13.88	165	579	2.934	ng	# 69
53) 3-Nitroaniline	14.23	138	19	3.275	ng	# 3
54) 2,4-Dinitrophenol	14.40	184	27	8.061	ng	# 1
57) 2,4-Dinitrotoluene	14.71	165	185	3.814	ng	# 68
62) 4-Nitroaniline	15.41	138	11	2.417	ng	# 1
65) 4,6-Dinitro-2-methylphenol	15.45	198	15	6.902	ng	# 1
80) Butylbenzylphthalate	20.43	149	157	3.061	ng	91
84) Bis(2-ethylhexyl)phthalate	20.96	149	240916	3.413	ng	# 52

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

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