

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061121\
 Data File : BP005876.D
 Acq On : 11 Jun 2021 18:04
 Operator : CG/JU
 Sample : M2613-21 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B12(114-118)

Quant Time: Jun 11 18:32:44 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 13:53:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	1628	20.00	ng	# 0.00
21) Naphthalene-d8	10.757	136	5550	20.00	ng	# 0.00
39) Acenaphthene-d10	14.581	164	2449	20.00	ng	0.00
64) Phenanthrene-d10	17.339	188	6225	20.00	ng	# 0.02
76) Chrysene-d12	21.392	240	10321	20.00	ng	# 0.00
86) Perylene-d12	23.810	264	13829	20.00	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	1569	15.46	ng	0.02
7) Phenol-d6	7.163	99	15362	116.82	ng	0.05
23) Nitrobenzene-d5	9.122	82	66	0.58	ng	0.01
42) 2,4,6-Tribromophenol	16.075	330	13	0.31	ng	0.01
45) 2-Fluorobiphenyl	13.251	172	855m	4.58	ng	0.05
79) Terphenyl-d14	19.875	244	5724	10.38	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.175	94	18796	143.08	ng	92
17) 2-Methylphenol	8.422	107	312	3.49	ng	# 69
20) 3+4-Methylphenols	8.775	107	1423	11.77	ng	# 76
75) Fluoranthene	19.333	202	1018	2.24	ng	78
77) Benzidine	19.533	184	5343	24.88	ng	# 96
81) Benzo(a)anthracene	21.380	228	1667	2.24	ng	94
83) Chrysene	21.433	228	1747	2.42	ng	# 85
87) Indeno(1,2,3-cd)pyrene	26.357	276	5168	4.37	ng	# 86
88) Benzo(b)fluoranthene	23.074	252	3755	3.96	ng	# 72
89) Benzo(k)fluoranthene	23.121	252	3294	3.59	ng	# 51
90) Benzo(a)pyrene	23.704	252	3111	3.49	ng	# 59
91) Dibenzo(a,h)anthracene	26.386	278	4921	4.83	ng	# 88
92) Benzo(g,h,i)perylene	27.133	276	7076	7.04	ng	# 84

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

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