

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072420\
 Data File : BF121010.D
 Acq On : 24 Jul 2020 18:49
 Operator : JU/CG
 Sample : L3360-01RX
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1ARX

Manual Integrations
 APPROVED

mohammad
 7/27/2020 12:19:33 PM

Quant Time: Jul 25 00:18:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	169140	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	607513	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	318997	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	553985	20.00	ng	0.00
76) Chrysene-d12	13.97	240	376853	20.00	ng	0.00
86) Perylene-d12	15.42	264	454288	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	370914	35.95	ng	0.00
7) Phenol-d6	6.44	99	858751	64.43	ng	0.00
23) Nitrobenzene-d5	7.37	82	640894	61.04	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	70986	20.33	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1219018	57.06	ng	0.00
79) Terphenyl-d14	12.92	244	1015660	58.59	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	238356	10.357	ng	99
71) Phenanthrene	11.36	178	657549	23.079	ng	98
72) Anthracene	11.40	178	169427	5.922	ng	98
73) Carbazole	11.56	167	111938	3.943	ng	99
75) Fluoranthene	12.54	202	1405890	44.517	ng	99
78) Pyrene	12.77	202	1247464	46.820	ng	99
81) Benzo(a)anthracene	13.96	228	691145	30.286	ng	98
83) Chrysene	14.00	228	660533	27.543	ng	97
87) Indeno(1,2,3-cd)pyrene	16.84	276	375315m	11.879	ng	
88) Benzo(b)fluoranthene	15.00	252	811847m	29.565	ng	
89) Benzo(k)fluoranthene	15.02	252	273423m	10.918	ng	
90) Benzo(a)pyrene	15.36	252	592682	23.657	ng	97
91) Dibenzo(a,h)anthracene	16.85	278	102946	3.989	ng	# 89
92) Benzo(g,h,i)perylene	17.27	276	359593	14.132	ng	99

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

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