

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010620\
 Data File : BP001541.D
 Acq On : 06 Jan 2020 21:35
 Operator : JU
 Sample : L1010-01 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-01-010220

Manual Integrations
 APPROVED

mohammad
 1/7/2020 12:31:41 PM

Quant Time: Jan 07 03:40:01 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 06 16:38:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	55884	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	214021	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	141094	20.00	ng	0.00
64) Phenanthrene-d10	17.08	188	305654	20.00	ng	0.00
76) Chrysene-d12	21.18	240	247634	20.00	ng	0.00
87) Perylene-d12	23.43	264	264203	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	128099	38.51	ng	0.00
7) Phenol-d6	6.88	99	188619	42.10	ng	0.00
23) Nitrobenzene-d5	8.83	82	126380	29.17	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	76317	47.79	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	281935	29.38	ng	0.00
79) Terphenyl-d14	19.66	244	398239	32.05	ng	0.00
Target Compounds						
71) Phenanthrene	17.12	178	90205	5.441	ng	98
75) Fluoranthene	19.10	202	162310	8.088	ng	100
78) Pyrene	19.45	202	163867	9.119	ng	98
81) Benzo(a)anthracene	21.16	228	65828	3.789	ng	92
83) Chrysene	21.22	228	65341m	3.922	ng	
88) Benzo(b)fluoranthene	22.76	252	74903m	4.579	ng	
90) Benzo(a)pyrene	23.33	252	53549	3.497	ng	# 94
92) Benzo(g,h,i)perylene	26.39	276	36657	2.415	ng	# 93

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

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