

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
 Data File : BP001663.D
 Acq On : 17 Jan 2020 18:27
 Operator : JU
 Sample : L1010-06 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-02-011620

Manual Integrations
 APPROVED

mohammad
 1/21/2020 7:53:04 AM

Quant Time: Jan 18 04:02:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 21:52:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	58966	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	204964	20.00	ng	-0.01
39) Acenaphthene-d10	14.28	164	110977	20.00	ng	0.00
64) Phenanthrene-d10	17.04	188	210647	20.00	ng	-0.01
76) Chrysene-d12	21.15	240	212287	20.00	ng	-0.01
87) Perylene-d12	23.38	264	226244	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	156086	51.85	ng	0.00
7) Phenol-d6	6.83	99	226625	47.11	ng	0.00
23) Nitrobenzene-d5	8.78	82	153980	32.83	ng	-0.01
42) 2,4,6-Tribromophenol	15.78	330	51983	30.77	ng	-0.01
45) 2-Fluorobiphenyl	12.89	172	260887	34.56	ng	-0.01
79) Terphenyl-d14	19.63	244	309047	26.63	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	10.46	128	28591	2.643	ng	95
52) Acenaphthene	14.35	154	15098	2.392	ng	97
58) Fluorene	15.34	166	25452	2.946	ng	95
71) Phenanthrene	17.08	178	199565	17.421	ng	99
72) Anthracene	17.17	178	39529	3.545	ng	97
75) Fluoranthene	19.06	202	212034	14.067	ng	99
78) Pyrene	19.42	202	214588	13.761	ng	97
81) Benzo(a)anthracene	21.13	228	94201	6.372	ng	96
83) Chrysene	21.18	228	85382	5.926	ng	97
86) Indeno(1,2,3-cd)pyrene	25.63	276	35140	2.084	ng	99
88) Benzo(b)fluoranthene	22.72	252	81623m	5.725	ng	
89) Benzo(k)fluoranthene	22.75	252	34141m	2.456	ng	
90) Benzo(a)pyrene	23.27	252	65372	4.829	ng	# 97
92) Benzo(a,h,i)perylene	26.31	276	33295	2.398	ng	# 93

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

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