

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010320\
 Data File : BP001528.D
 Acq On : 04 Jan 2020 02:50
 Operator : JU
 Sample : L1011-08 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MC-010220-2-5-COMP-B7

Manual Integrations
 APPROVED

mohammad
 1/6/2020 11:43:11 AM

Quant Time: Jan 04 04:04:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 03 15:57:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	63861	20.00	ng	-0.01
21) Naphthalene-d8	10.47	136	232761	20.00	ng	-0.01
39) Acenaphthene-d10	14.33	164	147138	20.00	ng	-0.01
64) Phenanthrene-d10	17.09	188	326795	20.00	ng	-0.01
76) Chrysene-d12	21.19	240	295355	20.00	ng	-0.01
87) Perylene-d12	23.44	264	329884	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	60513	15.92	ng	0.02
7) Phenol-d6	6.90	99	86285	16.85	ng	0.02
23) Nitrobenzene-d5	8.84	82	55528	11.78	ng	0.00
42) 2,4,6-Tribromophenol	15.83	330	27291	16.39	ng	-0.01
45) 2-Fluorobiphenyl	12.95	172	119050	11.90	ng	-0.02
79) Terphenyl-d14	19.67	244	189305	12.77	ng	-0.01
Target Compounds						
31) Naphthalene	10.52	128	51698	4.223	ng	Qvalue 99
49) Acenaphthylene	14.05	152	57213	4.130	ng	99
52) Acenaphthene	14.39	154	75855	8.725	ng	97
55) Dibenzofuran	14.73	168	84186	6.322	ng	98
58) Fluorene	15.39	166	103178	10.172	ng	94
71) Phenanthrene	17.13	178	1333147	75.214	ng	98
72) Anthracene	17.22	178	371373	21.140	ng	99
73) Carbazole	17.50	167	119608	7.358	ng	100
75) Fluoranthene	19.12	202	2113779	98.514	ng	96
78) Pyrene	19.47	202	2030000	94.716	ng	99
81) Benzo(a)anthracene	21.17	228	867232	41.847	ng	95
83) Chrysene	21.23	228	785497	39.530	ng	97
86) Indeno(1,2,3-cd)pyrene	25.73	276	546834	22.611	ng	# 92
88) Benzo(b)fluoranthene	22.77	252	1068071	52.290	ng	# 96
89) Benzo(k)fluoranthene	22.81	252	289438m	14.793	ng	
90) Benzo(a)pyrene	23.35	252	846045m	44.249	ng	
91) Dibenzo(a,h)anthracene	25.74	278	124338	6.543	ng	# 81
92) Benzo(g,h,i)perylene	26.43	276	555388	29.303	ng	# 95

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

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